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**Groundwater Monitoring Report No. 1
(First Quarterly Sampling Event)**

for the

**Sheridan Disposal Services Superfund Site
Operable Unit 2
Waller County, Texas**

Prepared by

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1.0 INTRODUCTION

1.1 *Purpose and Objectives*

This First Quarter Groundwater Monitoring Report for the Sheridan Disposal Services Superfund Site Ground Water Migration Management Operable Unit 2 (OU2) has been prepared on behalf of the Sheridan Site Trust (SST) in accordance with the Record of Decision (ROD) signed September 22, 1989, the Statement of Work (SOW), the Ground Water Consent Decree (CD), except as modified by later agreement between SST and the U.S. Environmental Protection Agency (EPA), and the Groundwater Migration Management Workplan approved July 18, 2006.

The purpose of this report is to summarize the data collected during the first quarter groundwater monitoring event conducted on August 2-3, 2006.

1.2 *Site Location and Description*

The Sheridan Disposal Services Superfund Site is located in northern Waller County, Texas, approximately nine miles north-northwest of the City of Hempstead, Texas and two miles northwest of the intersection of Clark Bottom Road and Farm Road 1736. The property is bounded on the east, south and west sides by farm and ranch lands and on the north by the Brazos River. The site lies within the Gulf Coastal Plain Physiographic Province and is transitionally positioned between the Post Oak Savannah and Blackland Prairie Natural Regions of Texas.

The Site formerly encompassed approximately 110 acres and included a 42-acre evaporation system, a 12-acre lagoon, a 17-acre dike surrounding the former lagoon, and miscellaneous processing equipment.

1.3 *Operable Unit 2 History*

In the final closure plan submitted to the state by SDS, the Sheridan Disposal Services Superfund Site was considered one unit. It was not until the U.S. EPA was involved with the site that 2 operable units were established. The Source Control unit was designated OU1 and the Ground Water Migration Management unit was designated OU2.

The ROD for OU2 was signed by U.S. EPA on September 27, 1989. The 1989 ROD identified natural attenuation as the selected remedy. The Ground Water Migration Management Consent Decree, ROD and Statement of Work were lodged in federal court in December 1991, but weren't entered until October 22, 1997. The beginning of remedial action for OU2 was predicated on the completion of the remedial action for OU1 based on the assumption that without the source (sludge) available, the ground water should be cleaned by natural attenuation from biological activity, sorption and filtration.

1.4 *Operable Unit 2 Remedy*

The major components of the remedy for Sheridan OU2 include:

- Natural attenuation of the ground water;
- Monitoring of ground water to ensure that the ACL are not exceeded;

- Sampling and analysis of the Brazos River immediately downgradient and upgradient of the point of entry of ground water from the site to the river; and
- Development of a corrective action plan to ensure that protective levels are met at the point of potential exposure if the ACLs are exceeded.

2.0 ASSESSMENT MONITORING PROGRAM

2.1 *Record of Decision Requirements*

U.S. EPA has selected ACLs that are the appropriate ground water standard for the site as long as the conditions set forth below remain valid. ACLs are ground water protection standards that are used to assure that hazardous constituents found in the ground water do not pose a risk to human health or the environment. To ensure that the ACLs remain protective, the following conditions must continue to be met at the site:

- 1) The Brazos River must remain the discharge point for ground water from the site.
- 2) The Brazos River cannot be adversely impacted by the discharge of contaminated ground water into the river. To ensure that future adverse impacts from the site do not occur at the point of exposure for environmental receptors in the river, river water will be sampled to ensure that there is no statistically significant increase in contamination, as compared to upgradient locations.
- 3) The ground water use restrictions must be implemented and continued to ensure that affected ground water is not consumed and the integrity of the Brazos River as a hydraulic barrier to ground water flow is maintained. Groundwater restrictions specified in the Record of Decision and Consent Decree include: no groundwater use within 100 feet from the edge of the plume and the owner will take no action at the site without getting consent from EPA, including sale of site.

2.2 *Remedy Assessment Criteria*

Natural attenuation was chosen as the final remedy for groundwater. As part of the remedy selection process, ACLs were established for the groundwater protection standard. The ACL values were calculated by determining the volume of affected water entering the river at any time and factoring in the dilution which would occur in the river at historical low flow conditions.

COMPOUND	ALTERNATE CONCENTRATION LIMITS (mg/l)
Benzene	26
Tetrachloroethylene	41
Trans-1,2-Dichloroethylene	26
Trichloroethylene	26
Arsenic	260

The point of compliance for meeting the ACLs is the location where the ACLs must be met and is also the well location where ACLs are monitored. At the point of compliance, ACLs ensure that human health and the environment are protected at the point of exposure and no statistically significant increase in contamination occurs in the river.

3.0 SAMPLING AND ANALYSIS PROCEDURES

3.1 *Pre-Sampling Activities*

Prior to the start of groundwater and surface water sampling, the existing monitor wells, MW-6, MW-31, MW-34, MW-35, MW-37, and MW-39, were located in the field and evaluated for adequacy. These monitoring wells were surveyed for top-of-casing (TOC) elevations and the elevations were tied to the permanent survey monuments that were established during the remediation of OU1. Elevations were measured to an accuracy of 0.01 feet and were recorded relative to the USC and GS 1983 North American datum.

The monitoring wells were also visually inspected to check the integrity of the protective steel casing, the concrete pad, the polyvinyl chloride (PVC) riser pipe, and the total depth of the well. The existing concrete pad associated with monitoring well MW-31 was removed and replaced with a 2-foot by 2-foot x 6 inch thick concrete pad that was sloped to provide water drainage away from the monitoring well. New 2-inch PVC caps were added to monitoring wells MW-6, MW-35 and MW-37 and the protective steel casings associated with monitoring wells MW-6, MW-31, MW-34, MW-35, MW-37, and MW-39 were primed and painted with rust-resistant yellow paint.

The total depth of the monitoring well and the depth to groundwater in each monitoring well were measured. Because the monitoring wells appeared to be "silted in", each monitoring well was redeveloped prior to sampling. Redevelopment was accomplished through pumping until the pH, specific conductance and water clarity stabilized. Approximately 30 to 55 gallons of development water were removed from each monitoring well. The water is temporarily stored in 55-gallon drums located at each monitoring well pending disposal. Ground water sampling activities were not initiated until a minimum of 24 hours had elapsed since redevelopment.

3.2 *Ground Water Sampling*

Groundwater sampling for the constituents of concern was used to determine the presence and concentration of the constituents, and if ACLs were approached or exceeded. The measurement of water levels at the site was used to determine the ground water flow direction and gradient to ensure that the Brazos River is the receptor of ground water from the site. Sampling of water from the Brazos River ensured that there was no impact on the river from the ground water.

3.2.1 Sampling Procedures

Ground water samples were collected from each monitoring well using low flow sampling techniques to minimize the effects of sediment entrained in the sample during analysis. The methods described in the U.S. EPA guidance document titled "Low-Flow (Minimal Drawdown) Groundwater Sampling Procedures" by Puls & Barcelona (EPA/540/S-95/504) were followed as described in the following paragraphs.

A variable flow submersible pump intake was placed at the middle, or slightly above the middle, of the screened interval and a low flow rate was used to draw formation water through the screen and up to the tubing outport. The flow rate was on the order of 0.1 – 0.5 L/min to minimize stress (drawdown of the water in the well casing), thereby minimizing any potential for overlying and underlying stagnant water to

enter the pump intake. An in-line flow through cell was attached to the outport which allowed for a continual read-out of water quality parameters (i.e. pH, specific conductivity, temperature, dissolved oxygen, and Eh). Once these parameters had stabilized (indicative of formation water), the well was sampled regardless of the volume of water purged. Turbidity was also measured with intermittent samples using a HACH meter not attached to the flow through cell. Well purging operations during the sampling event were conducted with a YSI Water Quality Meter equipped with a flow through cell. All readings were recorded in the field logbook.

Upon the completion of sampling, the sample containers were labeled and placed on ice in laboratory supplied ice chests. The samples were shipped to the analytical laboratory at the completion of sampling with the proper chain-of-custody forms using an overnight delivery service. In addition to the ground water samples, a quality control sample consisting of one trip blank was also collected during the sampling event.

3.2.2 Analytical Methods

Samples were analyzed for volatile organic compounds (VOCs), semi-volatile organic compounds (SVOCs), pesticides, PCBs, and metals. The VOC analysis was performed using EPA SW-846 Method 8260B, SVOC analysis by EPA SW-846 8270C, pesticides by EPA SW-846 8081A, and PCBs by EPA SW-846 8082. Samples for metals analysis were filtered in the field with a 0.45 micron filter and submitted for analysis by EPA SW-846 6020/7470A. The specific constituents of concern included the following:

Volatile Organic Compounds		
Acetone	Chloroform	4-Methyl-2-pentanone (MIBK)
Benzene	Chloromethane	Styrene
Bromodichloromethane	1,1-Dichloroethane	1,1,2,2-Tetrachloroethane
Bromoform	1,1-Dichloroethene	Tetrachloroethene
Bromomethane	trans-1,2-Dichloroethene	Toluene
2-Butanone (MEK)	1,2-Dichloropropane	1,1,1-Trichloroethane
Carbon disulfide	cis-1,3-Dichloropropene	1,1,2-Trichloroethane
Carbon tetrachloride	trans-1,3-Dichloropropene	Trichloroethene
Chlorodibromomethane	Ethylbenzene	Vinyl acetate
Chlorobenzene	2-Hexanone	Vinyl chloride
Chloroethane	Methylene chloride	Xylenes

Semivolatile Organic Compounds		
Acenaphthene	m-Cresol	Hexachloroethane
Acenaphthylene	p-Cresol	Indeno(1,2,3-cd) pyrene
Anthracene	Di-n-butylphthalate	2-Methylnaphthalene
Benz(a)anthracene	Dibenz (a,h) anthracene	Naphthalene
Benzo(b)fluoranthene	1,2-Dichlorobenzene	2-Nitroaniline
Benzo(k)fluoranthene	1,3-Dichlorobenzene	3-Nitroaniline
Benzo(g,h,i)perylene	1,4-Dichlorobenzene	4-Nitroaniline
Benzo(a)pyrene	3,3'-Dichlorobenzidine	Nitrobenzene
Benzoic acid	2,4-Dichlorophenol	2-Nitrophenol
Benzyl alcohol	Diethyl phthalate	4-Nitrophenol
Bis(2-chloroethoxy) methane	2,4-Dimethylphenol	N-Nitrosodimethylamine
Bis(2-chloroethyl) ether	Dimethylphthalate	N-Nitrosodiphenylamine
Bis(2-chloroisopropyl) ether	4,6-Dinitro-2-methylphenol	N-Nitrosodi-n-propylamine
Bis(2-ethylhexyl) phthalate	2,4-Dinitrophenol	Pentachlorophenol
4-Bromophenyl phenyl ether	2,4-Dinitrotoluene	Phenanthrene
Butyl benzyl phthalate	2,6-Dinitrotoluene	Phenol
p-Chloroaniline	Di-n-octylphthalate	Pyrene
p-Chloro-m-cresol	Fluoranthene	1,2,4-Trichlorobenzene
2-Chloronaphthalene	Fluorene	2,4,5-Trichlorophenol
2-Chlorophenol	Hexachlorobenzene	2,4,6-Trichlorophenol
4-Chlorophenyl phenyl ether	Hexachlorobutadiene	
Chrysene	Hexachlorocyclopentadiene	

Metals		
Arsenic	Chromium	Selenium
Barium	Lead	Silver
Cadmium	Mercury	Zinc
	Nickel	

Pesticides/PCBs		
Aldrin	Dieldrin	Aroclor 1242
alpha-BHC	Endosulfan I	Aroclor 1254
Beta-BHC	Endosulfan II	Aroclor 1221
delta-BHC	Endosulfan sulfate	Aroclor 1232
gamma-BHC (Lindane)	Endrin	Aroclor 1248
Chlordane	Endrin ketone	Aroclor 1260
4,4'-DDT	Heptachlor	Aroclor 1216
4,4'-DDE	Heptachlor epoxide	Toxaphene
4,4'-DDD	Methoxychlor	

3.3 Surface Water Sampling

Surface water samples were collected from two locations in the Brazos River to ensure there is no impact to the river from the site. One sample point was adjacent to the point of projected horizontal and vertical entry of the plume into the river (considered the upstream sample for the purposes of the statistical evaluation), and the other was downstream of the site. (The next sampling event will be performed at locations upstream and adjacent to the entry point of the ground water from the site into the Brazos River). The current samples were collected in quadruplicate to provide an adequate database to perform statistical analysis.

Surface water sampling took place in conjunction with the ground water sampling.

3.3.1 Sampling Procedures

Sampling of the surface water took place from a boat launched into the river. Since the water depth at the sampling points was greater than 0.46 m (1.5 ft), the samples were collected at a depth of approximately 0.3 m (1 ft) below the water surface. A properly decontaminated Kemmerer bottle was used to collect the surface water samples. The sampling device was lowered to the predetermined depth in the water column so that the sampling end pieces (upper and lower stoppers) were pulled away from the sampling tube (body), allowing the water to be sampled to pass through this tube. When the Kemmerer bottle was at the required depth, the sampling device was closed. The sampler was then retrieved and the first 10 to 20 ml of sample was discharged to clear any potential contamination of the valve. The water sample was then transferred to the appropriate laboratory-supplied sample container. Those samples that were analyzed for metals were field filtered using a 0.45 micron filter prior to placement in the sample bottle.

Upon the completion of quadruplicate sampling, the sample containers were labeled and placed on ice in laboratory supplied ice chests. The samples were shipped to the analytical laboratory at the completion of sampling with the proper chain-of-custody forms using an overnight delivery service.

3.3.2 Analytical Methods

Samples were analyzed for VOCs, SVOCs, pesticides, PCBs, and metals. The VOC analysis was performed using EPA SW-846 Method 8260B, SVOC analysis by EPA SW-846 8270C, pesticides by EPA SW-846 8081A, and PCBs by EPA SW-846 8082. Samples for metals analysis were filtered in the field with a 0.45 micron filter and submitted for analysis by EPA SW-846 6010B/7470A. The specific constituents of concern included those presented in Section 3.2.2 of this report.

4.0 EVALUATION OF MONITORING DATA

4.1 Analytical Results

The groundwater and surface water monitoring results from the August 2-3, 2006 monitoring event are presented in Table 1. Data for constituents detected below reporting limits and qualified as estimated ("J") and constituents detected in the blank samples (B) were excluded from further evaluation. The laboratory analytical reports showing all detection limits, estimated J values and the blank results are included in Appendix A. As shown in Table 1.0, concentrations of benzene, tetrachloroethylene, trans-1,2-dichloroethylene, trichloroethylene, and arsenic are below the established ACLs.

Vinyl chloride was detected at a concentration of 0.011 mg/L in monitoring well MW-37 only. This concentration exceeds the Maximum Contaminant Level (MCL) of 0.002 mg/L, however it was not detected in the surface water sample. The concentration of vinyl chloride will be tracked closely, and if necessary, an ACL will be calculated using the same methodology as was used to determine the other ACLs.

With regard to a comparison of the August 2006 sampling results to the analytical results for the previous sampling event conducted in October 1987, the following conclusions can be drawn:

- Constituent concentrations in the groundwater collected from monitoring wells MW-6, MW-31, MW-35, and MW-39 generally remain unchanged. While the laboratory detection limits for the constituents of concern have become more precise, the constituents detected are within the same order of magnitude. This is illustrated graphically in Figure 2

4.2 Groundwater Gradient

The groundwater gradient and flow direction for the site were determined using the groundwater elevation data collected from the monitoring wells during the August 2-3, 2006 sampling event. These data are included in Table 2 and depicted on Figure 1. Based on the data collected during the sampling event, the groundwater flow direction is to the northeast towards the Brazos River, as it has historically been.

4.3 Statistical Analysis of Surface Water Sampling Data

A limited statistical analysis per Section 3.2 of the Statement of Work for OU2 was performed to compare the adjacent and downstream constituent concentrations for the surface water samples collected during the August 2-3, 2006 sampling event. Because no detectable concentrations of benzene, tetrachloroethylene, trans-1,2-dichloroethylene, and trichloroethylene were present in both the adjacent and downstream samples, a statistical analysis was not performed for these constituents beyond that of the sample mean. The background mean for each of these constituents was considered to be equal to the method detection limit of 0.0002 mg/l for benzene, PCE and trans-1,2-DCE and 0.00032 mg/l for TCE. Since no constituents were detected above these method detection limits in the adjacent or downstream samples, it stands that the background mean was not exceeded in the downstream samples.

Arsenic concentrations, however, were detected slightly above the detection limit in both the adjacent and downstream surface water samples. Therefore, the sample mean for both the adjacent and downstream samples was calculated. The results are as follows:

- Adjacent sample mean: 0.004675
- Downstream sample mean: 0.00495

The average arsenic concentration for the downstream sample slightly exceeds the average arsenic concentration for the adjacent sample (considered the upstream sample for the purposes of the statistical evaluation). For this reason, the Dunnett's test was performed to determine if a statistically significant increase in the concentration of arsenic has occurred. Based on the calculations, the downstream samples do not have arsenic levels that are significantly higher than the background sample. No statistical difference was found between the upstream and downstream average sample concentrations. Detailed calculations are attached as Appendix B.

4.4 Further Action

The concentrations of the constituents of concern in the groundwater or surface water did not exceed the established trigger levels for increased monitoring, as presented below.

Trigger Levels for Increased Frequency Of Groundwater Monitoring	
COMPOUND	TRIGGER LEVEL (mg/L)
Benzene	1
Tetrachloroethylene	2
Trans-1,2-Dichloroethylene	1
Trichloroethylene	1
Arsenic	10

Therefore, based on the results from the August 2-3, 2006 sampling event, no further action with respect to an increase in the monitoring frequency is required.

5.0 REFERENCES

ERM-Southwest, Inc. 1990. Statement of Work for Remedial Design and Remedial Action, Ground Water Operable Unit. W.O. #91-21. Houston, Texas.

U.S. Environmental Protection Agency. 1997. Consent Decree, Civil Action No. H-91-3529. EPA Region VI, Dallas, Texas.

U.S. Environmental Protection Agency. 1989. Record of Decision for Sheridan Disposal Services Site. EPA Region VI, Dallas, Texas.

Table 1.0
 Groundwater and Surface Water Sampling Results

SHERIDAN DISPOSAL SERVICES SUPERFUND SITE
 GROUND WATER OPERABLE UNIT 2
 ANALYTICAL WATER RESULTS

Compound	Date	Benzene	Tetrachloroethylene	Trans-1,2-Dichloroethylene	Trichloroethylene	Total Arsenic	Vinyl Chloride
Alternate Concentration Limit		26	41	26	26	260	
Trigger for RAP Preparation		4	6	4	4	40	
Trigger for Increased Monitoring		1	2	1	1	10	
MW-6	08/03/06	BDL	BDL	BDL	BDL	ND	BDL
MW-31	08/03/06	BDL	BDL	BDL	BDL	0.0023	BDL
MW-34	08/03/06	0.067	BDL	0.0012	ND	0.0058	BDL
MW-35	08/03/06	ND	BDL	BDL	BDL	BDL	BDL
MW-37	08/03/06	ND	BDL	ND	ND	ND	0.011
MW-39	08/03/06	BDL	BDL	BDL	BDL	0.051	BDL
R1-A	08/02/06	BDL	BDL	BDL	BDL	0.0047	BDL
R1-B	08/02/06	BDL	BDL	BDL	BDL	ND	BDL
R1-C	08/02/06	BDL	BDL	BDL	BDL	ND	BDL
R1-D	08/02/06	BDL	BDL	BDL	BDL	0.0054	BDL
R2-A ¹	08/02/06	BDL	BDL	BDL	BDL	0.0051	BDL
R2-B	08/02/06	BDL	BDL	BDL	BDL	0.0058	BDL
R2-C	08/02/06	BDL	BDL	BDL	BDL	ND	BDL
R2-D	08/02/06	BDL	BDL	BDL	BDL	ND	BDL

Note - all concentrations in mg/L

1 - Upgradient Brazos River Sample

ND - estimated value below reporting limit

BDL - below detection limit

TABLE 2
SHERIDAN DISPOSAL SERVICES SUPERFUND SITE
GROUNDWATER OPERABLE UNIT 2
WELL DATA

Monitoring Well ID No.	Sample Date	Ground Elevation (ft amsl)	TOC Elevation (ft amsl)	Standpipe Stickup (+) Stickdown (-)	Total Well Depth (ft from gs)	Casing/ Screen Diameter (inches)	Screened Interval (ft from gs)	Depth to Water (ft from gs)	Depth to Water (ft from TOC)	Water Elevation (ft amsl)
MW-6	08/03/06	164.46	167.58	3.12	95.21	2	80-95	33.41	36.53	131.05
MW-31	08/03/06	166.70	168.67	1.97	65.01	4	25-60	35.34	37.31	131.36
MW-34	08/03/06	171.07	173.45	2.38	65.50	4	26-61	42.78	45.16	128.29
MW-35	08/03/06	171.32	173.39	2.07	105.02	2	80-100	41.44	43.51	129.88
MW-37	08/03/06	161.83	164.09	2.26	59.70	4	25-55	36.65	38.91	125.18
MW-39	08/03/06	164.81	166.41	1.60	59.00	4	34-54	34.15	35.75	130.66

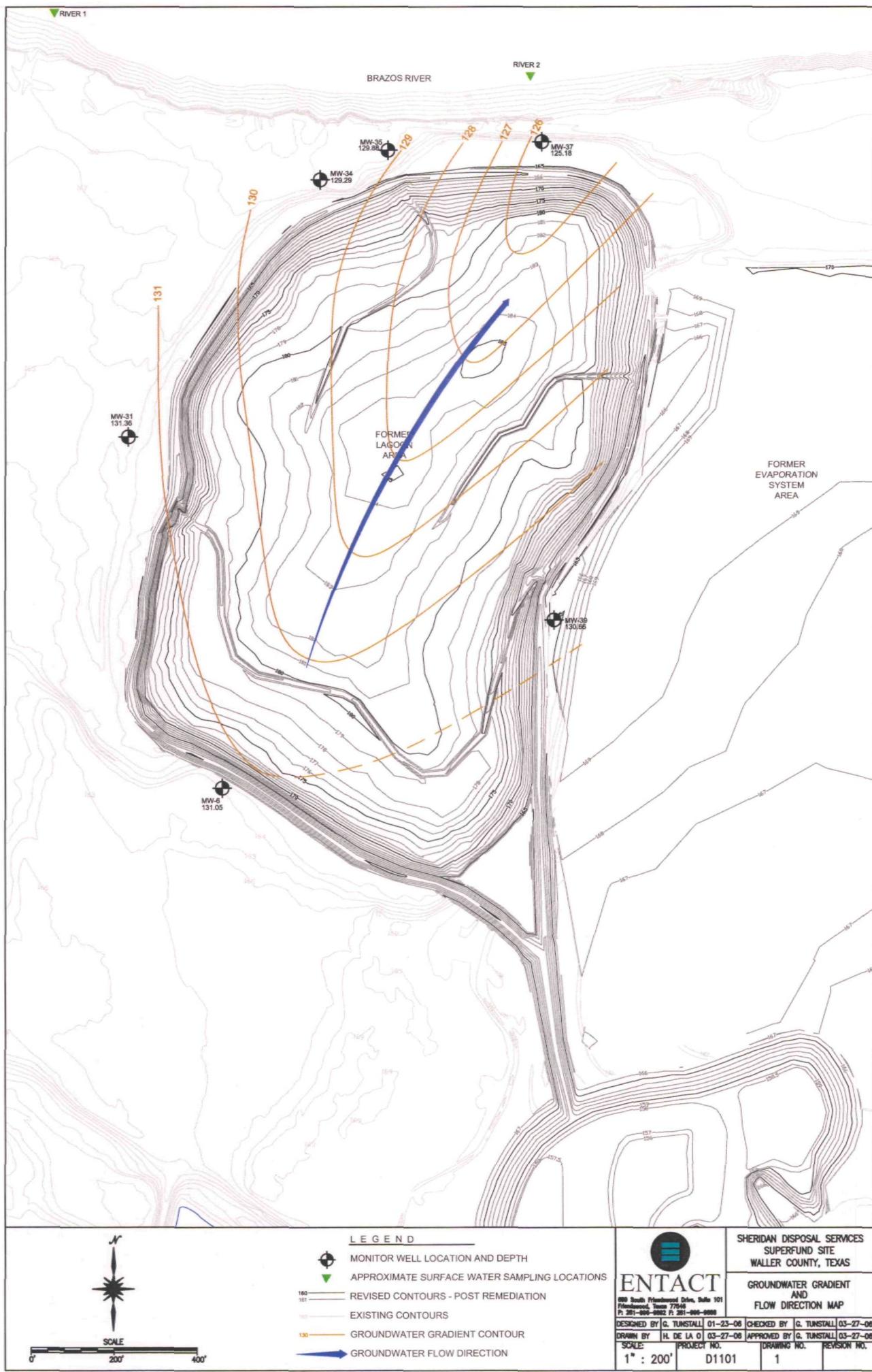
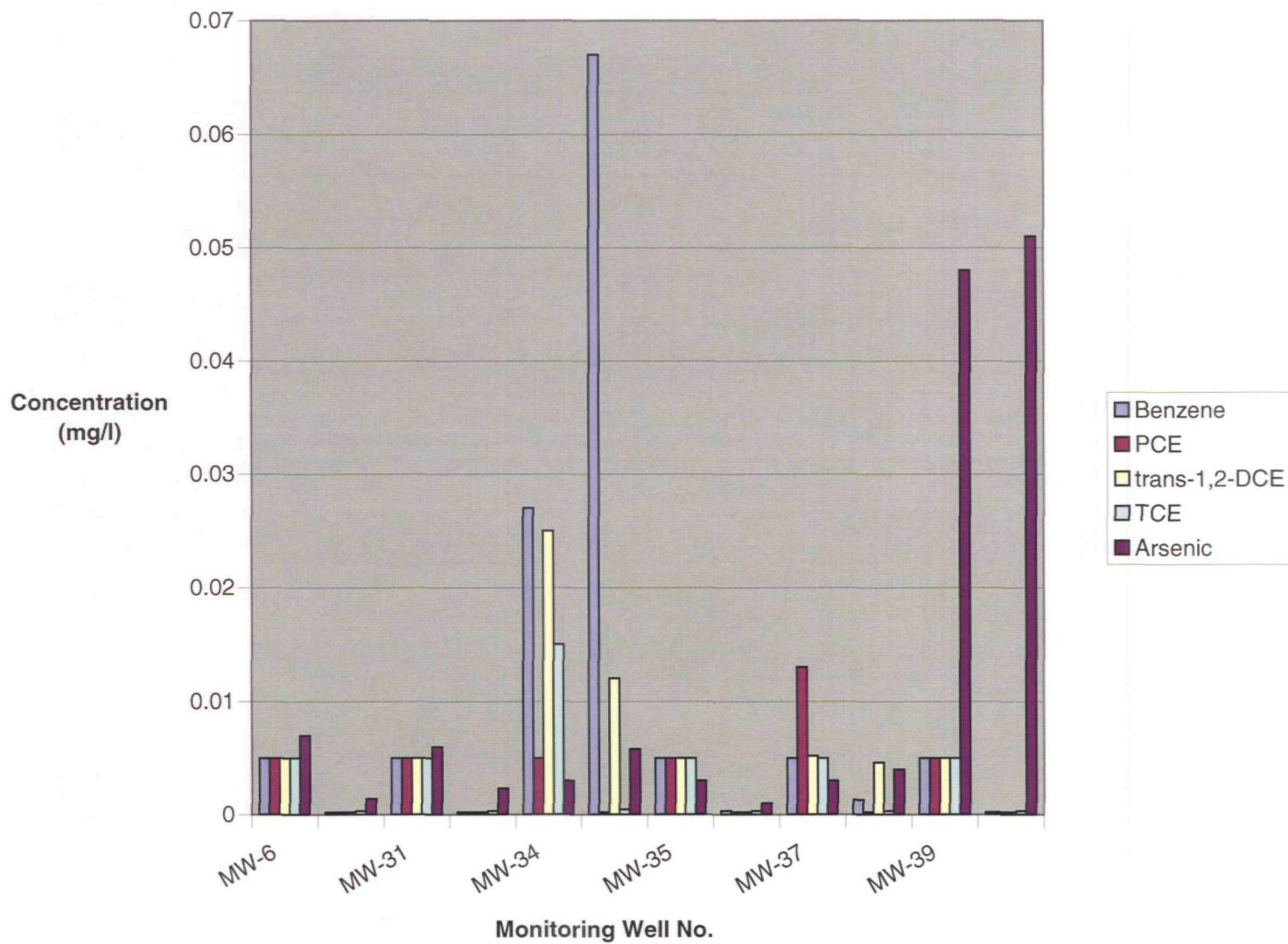


FIGURE 2
Graphical Comparison of the Constituent Concentrations in Groundwater



Note: October 1987 and August 2006 concentrations presented for each monitoring well.



ENTACT

Appendix

A

Appendix A



STL

ANALYTICAL REPORT

Job Number: 560-1385-1

Job Description: D1631 Sheridan Superfund

For:
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Attention: Ms. Liz Scaggs

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The test results entered in this report meet all NELAC requirements for accredited parameters. Any exceptions to NELAC requirements are noted in the report. Pursuant to NELAC, this report may not be reproduced except in full, and with written approval from the laboratory. STL Corpus Christi Certifications and Approvals: NELAC TX T104704210-06-TX, NELAC KS E-10362, NELAC LA 03034, Oklahoma 9968, USDA Soil Permit S-42935 Revised.

Case Narrative for job: 560-J1385-1

Client: Entact Environmental Services, LLC

Date: 09/06/2006

Total Metals Analysis

Samples 560-1385-1 through 14 were analyzed for total metals using EPA Method 6020. The percent recovery result for selenium in the LCS associated with these samples was above the acceptance criteria. Selenium was not detected in samples 1-12 and 14 but was detected at a low concentration in sample 13.

Dissolved Metals Analysis

Samples 560-1385-1 through 14 were analyzed for total metals using EPA Method 6020. The percent recovery results for all metals except silver and lead in the LCS associated with these samples were above the acceptance criteria. The results for barium were at least 3x the reporting limit. The other metals were detected at low concentrations in the various samples.

Organochlorine Pesticides Analysis

Samples 560-1385-3, 8 and 14 were analyzed for organochlorine pesticides using EPA Method 8081A. The percent recovery results for the matrix spike duplicate associated with sample 3 were below the acceptance criteria. The percent recovery results for tetrachloro-m-xylene surrogate in samples 8 and 14 were outside the acceptance criteria. The associated LCS was within acceptable limits and the data are therefore reported.

Polychlorinated Biphenyls (PCB) Analysis

Samples 560-1385-1 through 14 were analyzed for PCB using EPA Method 8082. The percent recovery result for Aroclor 1260 in the matrix spikes associated with sample 2 were below the acceptance criteria. The associated LCS was within acceptable limits and the data are therefore reported. The percent recovery results for decachlorobiphenyl surrogate in these samples were below the acceptance criteria but the results for tetrachloro-m-xylene surrogate were within acceptable limits. Per the method, the results are reportable with acceptable recoveries for one surrogate.

Semivolatile Organics Analysis

Sample 560-1385-1 was analyzed for semivolatile organics using EPA Method 8270C. The percent recovery results or RPD for 2,4-dinitrophenol, 4,6-dinitro-2-methylphenol, n-nitrosodiphenylamine and 3,3'-dichlorobenzidine in the matrix spike pair associated with this sample were outside the acceptance criteria. The associated LCS was within acceptable limits and the data are therefore reported.

Volatile Organics Analysis

Samples 560-1385-3 through 14 were analyzed for volatile organics using EPA Method 8260B. The percent recovery result for vinyl acetate in the LCS associated with these samples was above the acceptance criteria. Vinyl acetate was not detected in the samples and the data are therefore reported.

EXECUTIVE SUMMARY - Detections

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Sample ID	Client Sample ID		Reporting Limit	Units	Method
Analyte		Result / Qualifier			
560-1385-1	R2-A				
Acetone		3.6	J	100	ug/L
Toluene		0.33	J	5.0	ug/L
Xylenes, Total		0.44	J	15	ug/L
As		5.1		5.0	ug/L
Ba		190	B	50	ug/L
Ni		2.8	J B	5.0	ug/L
<i>Dissolved</i>					
As		6.3	*	5.0	6020
Ba		220	* B	50	6020
Cr		1.6	J *	20	6020
Ni		3.5	J * B	5.0	6020
Se		3.1	J *	5.0	6020
560-1385-2	R2-B				
Acetone		2.9	J	100	ug/L
Toluene		0.34	J	5.0	ug/L
Xylenes, Total		0.40	J	15	ug/L
gamma-BHC (Lindane)		0.0051	J	0.050	ug/L
As		5.8		5.0	6020
Ba		210	B	50	6020
Cr		1.4	J	20	6020
Ni		2.7	J B	5.0	6020
<i>Dissolved</i>					
As		5.8	*	5.0	6020
Ba		190	* B	50	6020
Cr		1.3	J *	20	6020
Ni		2.5	J * B	5.0	6020
Pb		4.5	J	5.0	6020
Se		4.5	J *	5.0	6020

EXECUTIVE SUMMARY - Detections

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Sample ID	Client Sample ID	Reporting				
Analyte		Result / Qualifier		Limit	Units	Method
560-1385-3	R2-C					
Acetone		4.0	J	100	ug/L	8260B
Toluene		0.22	J	5.0	ug/L	8260B
Xylenes, Total		0.39	J	15	ug/L	8260B
As		4.3	J	5.0	ug/L	6020
Ba		180	B	50	ug/L	6020
Ni		1.9	J B	5.0	ug/L	6020
Hg		0.00013	J	0.0020	mg/L	7470A
<i>Dissolved</i>						
As		4.4	J *	5.0	ug/L	6020
Ba		180	* B	50	ug/L	6020
Ni		1.4	J * B	5.0	ug/L	6020
560-1385-4	R2-D					
Acetone		4.1	J	100	ug/L	8260B
Toluene		0.21	J	5.0	ug/L	8260B
As		4.6	J	5.0	ug/L	6020
Ba		190	B	50	ug/L	6020
Cr		1.1	J	20	ug/L	6020
Ni		2.3	J B	5.0	ug/L	6020
Zn		53	J	100	ug/L	6020
<i>Dissolved</i>						
As		4.1	J *	5.0	ug/L	6020
Ba		180	* B	50	ug/L	6020
Ni		2.1	J * B	5.0	ug/L	6020
560-1385-5	R1-A					
Acetone		3.0	J	100	ug/L	8260B
Toluene		0.38	J	5.0	ug/L	8260B
Xylenes, Total		0.52	J	15	ug/L	8260B
As		4.7	J	5.0	ug/L	6020
Ba		180	B	50	ug/L	6020
Ni		2.0	J B	5.0	ug/L	6020
<i>Dissolved</i>						
As		4.8	J *	5.0	ug/L	6020
Ba		210	* B	50	ug/L	6020
Ni		2.3	J * B	5.0	ug/L	6020

EXECUTIVE SUMMARY - Detections

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Sample ID	Client Sample ID		Reporting Limit	Units	Method
Analyte		Result / Qualifier			
560-1385-6	R1-B				
Acetone		0.85	J	100	ug/L
Toluene		0.47	J	5.0	ug/L
Xylenes, Total		0.63	J	15	ug/L
As		4.5	J	5.0	ug/L
Ba		190	B	50	ug/L
Ni		1.8	J B	5.0	ug/L
<i>Dissolved</i>					
As		4.8	J *	5.0	ug/L
Ba		190	* B	50	ug/L
Ni		1.2	J * B	5.0	ug/L
560-1385-7	R1-C				
Acetone		2.0	J	100	ug/L
Toluene		0.27	J	5.0	ug/L
Xylenes, Total		0.38	J	15	ug/L
As		4.1	J	5.0	ug/L
Ba		170	B	50	ug/L
Ni		1.7	J B	5.0	ug/L
<i>Dissolved</i>					
As		4.0	J *	5.0	ug/L
Ba		180	* B	50	ug/L
Ni		1.3	J * B	5.0	ug/L
560-1385-8	R1-D				
Acetone		3.4	J	100	ug/L
Toluene		0.23	J	5.0	ug/L
Xylenes, Total		0.36	J	15	ug/L
As		5.4		5.0	ug/L
Ba		220	B	50	ug/L
Ni		2.3	J B	5.0	ug/L
<i>Dissolved</i>					
As		4.3	J *	5.0	ug/L
Ba		190	* B	50	ug/L
Ni		1.5	J * B	5.0	ug/L
Hg		0.00017	J	0.0020	mg/L

EXECUTIVE SUMMARY - Detections

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Sample ID	Client Sample ID		Reporting		
Analyte		Result / Qualifier	Limit	Units	Method
560-1385-9	MW-6				
Vinyl chloride		0.73	J	5.0	ug/L
As		1.4	J	5.0	ug/L
Ba		1200	B	50	ug/L
Cr		3.0	J	20	ug/L
Ni		2.1	J B	5.0	ug/L
Hg		0.00016	J	0.0020	mg/L
<i>Dissolved</i>					
Ba		1000	* B	50	ug/L
560-1385-10	MW-31				
Chloromethane		0.48	J	5.0	ug/L
Bis(2-ethylhexyl) phthalate		2.0	J	10	ug/L
As		23		5.0	ug/L
Ba		400	B	50	ug/L
Hg		0.00023	J	0.0020	mg/L
<i>Dissolved</i>					
As		11	*	5.0	ug/L
Ba		410	* B	50	ug/L
560-1385-11	MW-34				
Chloromethane		1.1	J	5.0	ug/L
Vinyl chloride		1.7	J	5.0	ug/L
1,1-Dichloroethene		0.91	J	5.0	ug/L
Acetone		22	J	100	ug/L
trans-1,2-Dichloroethene		12		5.0	ug/L
1,1-Dichloroethane		0.24	J	5.0	ug/L
Benzene		67		5.0	ug/L
Trichloroethene		0.44	J	5.0	ug/L
Toluene		0.23	J	5.0	ug/L
Chlorobenzene		0.52	J	5.0	ug/L
Methyl Ethyl Ketone		2.0	J	5.0	ug/L
As		5.8		5.0	ug/L
Ba		980	B	50	ug/L
<i>Dissolved</i>					
As		3.3	J *	5.0	ug/L
Ba		1100	* B	50	ug/L
Hg		0.00013	J	0.0020	mg/L

EXECUTIVE SUMMARY - Detections

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Sample ID	Client Sample ID	Reporting			
Analyte		Result / Qualifier	Limit	Units	Method
560-1385-12	MW-35				
Acetone		2.5	J	100	ug/L
Benzene		0.33	J	5.0	ug/L
Bis(2-ethylhexyl) phthalate		7.4	J	10	ug/L
Ba		150	B	50	ug/L
Ni		2.1	J B	5.0	ug/L
<i>Dissolved</i>					
Ba		160	* B	50	ug/L
560-1385-13	MW-37				
Chloromethane		0.44	J	5.0	ug/L
Vinyl chloride		11		5.0	ug/L
Carbon disulfide		0.22	J	5.0	ug/L
Acetone		4.7	J	100	ug/L
trans-1,2-Dichloroethene		4.6	J	5.0	ug/L
Benzene		1.3	J	5.0	ug/L
Trichloroethene		0.32	J	5.0	ug/L
Toluene		0.47	J	5.0	ug/L
Chlorobenzene		1.5	J	5.0	ug/L
Methyl Ethyl Ketone		0.91	J	5.0	ug/L
As		4.0	J	5.0	ug/L
Ba		490	B	50	ug/L
Ni		1.4	J B	5.0	ug/L
Se		6.2	*	5.0	ug/L
<i>Dissolved</i>					
As		3.4	J *	5.0	ug/L
Ba		450	* B	50	ug/L
Ni		2.5	J * B	5.0	ug/L
Se		5.3	*	5.0	ug/L

EXECUTIVE SUMMARY - Detections

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Sample ID	Client Sample ID	Reporting			
Analyte		Result / Qualifier	Limit	Units	Method
560-1385-14	MW-39				
Chloromethane		0.51	J	5.0	ug/L
Acetone		4.8	J	100	ug/L
Bis(2-ethylhexyl) phthalate		9.6	J	10	ug/L
Heptachlor		0.0081	J	0.050	ug/L
Endosulfan I		0.020	J	0.050	ug/L
As		51		5.0	ug/L
Ba		240	B	50	ug/L
Ni		12	B	5.0	ug/L
<i>Dissolved</i>					
As		51	*	5.0	6020
Ba		320	* B	50	6020
Ni		1.8	J * B	5.0	6020

METHOD SUMMARY

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Description	Lab Location	Method	Preparation Method
Matrix: Water			
Volatile Organic Compounds by GC/MS	STL CC	SW846 8260B	
Purge-and-Trap	STL CC		SW846 5030B
Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)	STL CC	SW846 8270C	
Continuous Liquid-Liquid Extraction	STL CC		SW846 3520C
Organochlorine Pesticides by Gas Chromatography	STL CC	SW846 8081A	
Continuous Liquid-Liquid Extraction/Shared Prep	STL CC		SW846 3520C
Polychlorinated Biphenyls (PCBs) by Gas Chromatography	STL CC	SW846 8082	
Continuous Liquid-Liquid Extraction/Shared Prep	STL CC		SW846 3520C
Inductively Coupled Plasma - Mass Spectrometry	STL CC	SW846 6020	
Acid Digestion of Aqueous Samples and Extracts	STL CC		SW846 3010A
Sample Filtration performed in the Field	STL CC		FIELD_FLTRD
Mercury in Liquid Waste (Manual Cold Vapor Technique)	STL CC	SW846 7470A	
Mercury in Liquid Waste (Manual Cold Vapor	STL CC		SW846 7470A
Sample Filtration performed in the Field	STL CC		FIELD_FLTRD

LAB REFERENCES:

STL CC = STL Corpus Christi

METHOD REFERENCES:

SW846 - "Test Methods For Evaluating Solid Waste, Physical/Chemical Methods", Third Edition, November 1986 And Its Updates.

METHOD / ANALYST SUMMARY

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method	Analyst	Analyst ID
SW846 8260B	Michalk, Kevin	KRM
SW846 8270C	Craig, Bronson	BC
SW846 8081A	Williams, Sharon	SEW
SW846 8082	Williams, Sharon	SEW
SW846 6020	Mathewson, John E	JEM
SW846 7470A	Theriault, Ray	RT

SAMPLE SUMMARY

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Sample ID	Client Sample ID	Client Matrix	Date/Time Sampled	Date/Time Received
560-1385-1	R2-A	Water	08/02/2006 1123	08/04/2006 1040
560-1385-2	R2-B	Water	08/02/2006 1147	08/04/2006 1040
560-1385-3	R2-C	Water	08/02/2006 1215	08/04/2006 1040
560-1385-4	R2-D	Water	08/02/2006 1241	08/04/2006 1040
560-1385-5	R1-A	Water	08/02/2006 1621	08/04/2006 1040
560-1385-6	R1-B	Water	08/02/2006 1643	08/04/2006 1040
560-1385-7	R1-C	Water	08/02/2006 1709	08/04/2006 1040
560-1385-8	R1-D	Water	08/02/2006 1729	08/04/2006 1040
560-1385-9	MW-6	Water	08/03/2006 0759	08/04/2006 1040
560-1385-10	MW-31	Water	08/03/2006 0920	08/04/2006 1040
560-1385-11	MW-34	Water	08/03/2006 1040	08/04/2006 1040
560-1385-12	MW-35	Water	08/03/2006 1155	08/04/2006 1040
560-1385-13	MW-37	Water	08/03/2006 1253	08/04/2006 1040
560-1385-14	MW-39	Water	08/03/2006 1358	08/04/2006 1040
560-1385-15TB	TRIP BLANK	Water	08/03/2006 0000	08/04/2006 1040

SAMPLE RESULTS

Ms. Liz Scaggs
 Entact Environmental Services, LLC
 3129 Bass Pro Drive
 Grapevine, TX 76051

Job Number: 560-1385-1

Client Sample ID: R2-A
 Lab Sample ID: 560-1385-1

Date Sampled: 08/02/2006 1123
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:		08/07/2006 1836			
Prep Method: 5030B	Date Prepared:		08/07/2006 1836			
Chloromethane	0.39	U	ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	3.6	J	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.33	J	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.44	J	ug/L	0.35	15	1.0
Surrogate					Acceptance Limits	
Dibromofluoromethane (Surr)	109		%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	113		%		70 - 120	
Toluene-d8 (Surr)	99		%		80 - 120	
4-Bromofluorobenzene (Surr)	92		%		75 - 120	

Ms. Liz Scaggs
Entact Environmental Services, LLC
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Grapevine, TX 76051

Job Number: 560-1385-1

Client Sample ID: R2-A
Lab Sample ID: 560-1385-1

Date Sampled: 08/02/2006 1123
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006	1449		
Prep Method: 3520C	Date Prepared:	08/08/2006	1230		
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	50
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Chloronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

Ms. Liz Scaggs
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 Grapevine, TX 76051

Job Number: 560-1385-1

Client Sample ID: R2-A
 Lab Sample ID: 560-1385-1

Date Sampled: 08/02/2006 1123
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8270C			Date Analyzed:	08/09/2006 1449		
Prep Method: 3520C			Date Prepared:	08/08/2006 1230		
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	U	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0

	Acceptance Limits		
Surrogate			
2-Fluorophenol	56	%	10 - 120
Phenol-d5	60	%	12 - 120
Nitrobenzene-d5	66	%	30 - 120
2-Fluorobiphenyl	59	%	26 - 120
2,4,6-Tribromophenol	63	%	25 - 120
Terphenyl-d14	49	%	10 - 120

Method: 8081A	Date Analyzed:	08/14/2006 1914
Prep Method: 3520C	Date Prepared:	08/08/2006 1700
alpha-BHC	0.0056	U
beta-BHC	0.0056	U
delta-BHC	0.0025	U
Heptachlor	0.0059	U
Aldrin	0.0025	U

Ms. Liz Scaggs
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 3129 Bass Pro Drive
 Grapevine, TX 76051

Job Number: 560-1385-1

Client Sample ID: R2-A
Lab Sample ID: 560-1385-1

Date Sampled: 08/02/2006 1123
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:		08/14/2006 1914			
Prep Method: 3520C	Date Prepared:		08/08/2006 1700			
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate						Acceptance Limits
Tetrachloro-m-xylene	60		%			57 - 127
DCB Decachlorobiphenyl	19		%			10 - 152
Method: 8082	Date Analyzed:		08/15/2006 1939			
Prep Method: 3520C	Date Prepared:		08/08/2006 1700			
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate						Acceptance Limits
Tetrachloro-m-xylene	86		%			25 - 140
DCB Decachlorobiphenyl	27	X	%			42 - 133
Method: 6020	Date Analyzed:		08/09/2006 0210			
Prep Method: 3010A	Date Prepared:		08/07/2006 1103			
Ag	1.0	U	ug/L	1.0	5.0	10
As	5.1	ug/L	1.0	5.0	10	
Ba	190	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R2-A

Lab Sample ID: 560-1385-1

Date Sampled: 08/02/2006 1123

Date Received: 08/04/2006 1040

Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0210			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	1.1	U ug/L	1.1	20	10
Ni	2.8	J B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0509			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	6.3	* ug/L	1.0	5.0	10
Ba	220	* B ug/L	1.0	50	10
Cd	1.0	U * ug/L	1.0	5.0	10
Cr	1.6	J * ug/L	1.1	20	10
Ni	3.5	J * B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	3.1	J * ug/L	1.0	5.0	10
Zn	50	U * ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/09/2006 1649			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/09/2006 1714			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-B
Lab Sample ID: 560-1385-2

Date Sampled: 08/02/2006 1147
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:		08/07/2006 1901			
Prep Method: 5030B	Date Prepared:		08/07/2006 1901			
Chloromethane	0.39	U	ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	2.9	J	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichlormethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.34	J	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.40	J	ug/L	0.35	15	1.0
Surrogate						Acceptance Limits
Dibromofluoromethane (Surr)	108		%			80 - 120
1,2-Dichloroethane-d4 (Surr)	113		%			70 - 120
Toluene-d8 (Surr)	101		%			80 - 120
4-Bromofluorobenzene (Surr)	94		%			75 - 120

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Job Number: 560-1385-1

Client Sample ID: R2-B
 Lab Sample ID: 560-1385-2

Date Sampled: 08/02/2006 1147
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1517			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	1.0
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Chloronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

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Job Number: 560-1385-1

Client Sample ID: R2-B
 Lab Sample ID: 560-1385-2

Date Sampled: 08/02/2006 1147
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1517			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	ug/L	0.65	10	1.0
Phenanthrene	0.50	ug/L	0.50	10	1.0
Anthracene	0.50	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	ug/L	0.50	10	1.0
Fluoranthene	0.50	ug/L	0.50	10	1.0
Pyrene	0.50	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	ug/L	0.50	10	1.0
Chrysene	0.50	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	ug/L	1.3	10	1.0
Benzoic acid	20	ug/L	20	50	1.0
Surrogate				Acceptance Limits	
2-Fluorophenol	62.	%		10 - 120	
Phenol-d5	66	%		12 - 120	
Nitrobenzene-d5	73	%		30 - 120	
2-Fluorobiphenyl	70	%		26 - 120	
2,4,6-Tribromophenol	68	%		25 - 120	
Terphenyl-d14	40	%		10 - 120	
Method: 8081A	Date Analyzed:	08/14/2006 1937			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
alpha-BHC	0.0056	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-B
 Lab Sample ID: 560-1385-2

Date Sampled: 08/02/2006 1147
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed: 08/14/2006 1937					
Prep Method: 3520C	Date Prepared: 08/08/2006 1700					
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0051	J	ug/L	0.0027	0.050	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	58		%		57 - 127	
DCB Decachlorobiphenyl	17		%		10 - 152	
Method: 8082	Date Analyzed: 08/15/2006 1956					
Prep Method: 3520C	Date Prepared: 08/08/2006 1700					
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	83		%		25 - 140	
DCB Decachlorobiphenyl	26	X	%		42 - 133	
Method: 6020	Date Analyzed: 08/09/2006 0216					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Ag	1.0	U	ug/L	1.0	5.0	10
As	5.8	ug/L	1.0	5.0	10	
Ba	210	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R2-B
Lab Sample ID: 560-1385-2

Date Sampled: 08/02/2006 1147
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0216				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Cr	1.4	J	ug/L	1.1	20	10
Ni	2.7	J B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U	ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0527				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	5.8	*	ug/L	1.0	5.0	10
Ba	190	* B	ug/L	1.0	50	10
Cd	1.0	U *	ug/L	1.0	5.0	10
Cr	1.3	J *	ug/L	1.1	20	10
Ni	2.5	J * B	ug/L	1.0	5.0	10
Pb	4.5	J	ug/L	1.0	5.0	10
Se	4.5	J *	ug/L	1.0	5.0	10
Zn	50	U *	ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/09/2006 1651				
Prep Method: 7470A	Date Prepared:	08/09/2006 1130				
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/09/2006 1717				
Prep Method: 7470A	Date Prepared:	08/09/2006 1130				
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-C
 Lab Sample ID: 560-1385-3

Date Sampled: 08/02/2006 1215
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:	08/09/2006 1616			
Prep Method: 5030B	Date Prepared:	08/09/2006 1616			
Chloromethane	0.39	U ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U ug/L	0.20	5.0	1.0
Bromomethane	0.39	U ug/L	0.39	5.0	1.0
Chloroethane	0.40	U ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U ug/L	0.53	50	1.0
Acetone	4.0	J ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U ug/L	0.20	5.0	1.0
Chloroform	0.20	U ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Benzene	0.20	U ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Toluene	0.22	J ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U ug/L	0.20	5.0	1.0
Bromoform	0.50	U ug/L	0.50	5.0	1.0
Styrene	0.50	U ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U ug/L	0.47	5.0	1.0
Xylenes, Total	0.39	J ug/L	0.35	15	1.0
Surrogate				Acceptance Limits	
Dibromofluoromethane (Surr)	101	%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	104	%		70 - 120	
Toluene-d8 (Surr)	95	%		80 - 120	
4-Bromofluorobenzene (Surr)	92	%		75 - 120	

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Job Number: 560-1385-1

Client Sample ID: R2-C
 Lab Sample ID: 560-1385-3

Date Sampled: 08/02/2006 1215
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte		Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C		Date Analyzed:	08/09/2006	1545		
Prep Method: 3520C		Date Prepared:	08/08/2006	1230		
Phenol	0.50	U	ug/L	0.50	10	1.0
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10	1.0
2-Chlorophenol	0.50	U	ug/L	0.50	10	1.0
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10	1.0
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10	1.0
Benzyl alcohol	1.4	U	ug/L	1.4	20	1.0
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10	1.0
2-Methylphenol	0.50	U	ug/L	0.50	10	1.0
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10	1.0
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10	1.0
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10	1.0
Hexachloroethane	0.58	U	ug/L	0.58	10	1.0
Nitrobenzene	0.50	U	ug/L	0.50	10	1.0
2-Nitrophenol	0.50	U	ug/L	0.50	10	1.0
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10	1.0
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10	1.0
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10	1.0
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10	1.0
Naphthalene	0.50	U	ug/L	0.50	10	1.0
4-Chloroaniline	0.50	U	ug/L	0.50	10	1.0
Hexachlorobutadiene	0.50	U	ug/L	0.50	10	1.0
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10	1.0
2-Methylnaphthalene	0.50	U	ug/L	0.50	10	1.0
Hexachlorocyclopentadiene	50	U	ug/L	50	50	1.0
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10	1.0
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10	1.0
2-Choronaphthalene	0.50	U	ug/L	0.50	10	1.0
2-Nitroaniline	0.50	U	ug/L	0.50	50	1.0
Dimethyl phthalate	0.55	U	ug/L	0.55	10	1.0
Acenaphthylene	0.50	U	ug/L	0.50	10	1.0
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10	1.0
3-Nitroaniline	1.8	U	ug/L	1.8	50	1.0
Acenaphthene	0.57	U	ug/L	0.57	10	1.0
2,4-Dinitrophenol	20	U	ug/L	20	50	1.0
4-Nitrophenol	10	U	ug/L	10	50	1.0
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10	1.0
Diethyl phthalate	0.52	U	ug/L	0.52	10	1.0
Fluorene	0.61	U	ug/L	0.61	10	1.0
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10	1.0
4-Nitroaniline	1.5	U	ug/L	1.5	50	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-C
 Lab Sample ID: 560-1385-3

Date Sampled: 08/02/2006 1215
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:		08/09/2006 1545			
Prep Method: 3520C	Date Prepared:		08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	U	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0
Surrogate				Acceptance Limits		
2-Fluorophenol	54		%	10 - 120		
Phenol-d5	58		%	12 - 120		
Nitrobenzene-d5	63		%	30 - 120		
2-Fluorobiphenyl	62		%	26 - 120		
2,4,6-Tribromophenol	58		%	25 - 120		
Terphenyl-d14	34		%	10 - 120		
Method: 8081A	Date Analyzed:		08/14/2006 2001			
Prep Method: 3520C	Date Prepared:		08/08/2006 1700			
alpha-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	U	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	U	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	U	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-C
Lab Sample ID: 560-1385-3

Date Sampled: 08/02/2006 1215
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2001				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	60		%		57 - 127	
DCB Decachlorobiphenyl	19		%		10 - 152	
Method: 8082	Date Analyzed:	08/15/2006 2047				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	82		%		25 - 140	
DCB Decachlorobiphenyl	27	X	%		42 - 133	
Method: 6020	Date Analyzed:	08/09/2006 0222				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	4.3	J	ug/L	1.0	5.0	10
Ba	180	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R2-C
Lab Sample ID: 560-1385-3

Date Sampled: 08/02/2006 1215
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0222			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	1.1	U ug/L	1.1	20	10
Ni	1.9	J B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0534			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	4.4	J * ug/L	1.0	5.0	10
Ba	180	* B ug/L	1.0	50	10
Cd	1.0	U * ug/L	1.0	5.0	10
Cr	1.1	U * ug/L	1.1	20	10
Ni	1.4	J * B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U * ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/09/2006 1702			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	J mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/09/2006 1720			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-D
 Lab Sample ID: 560-1385-4

Date Sampled: 08/02/2006 1241
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B			Date Analyzed:	08/09/2006 1758		
Prep Method: 5030B			Date Prepared:	08/09/2006 1758		
Chloromethane	0.39	U	ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	4.1	J	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U *	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.21	J	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U	ug/L	0.35	15	1.0
Surrogate						Acceptance Limits
Dibromofluoromethane (Surr)	103		%			80 - 120
1,2-Dichloroethane-d4 (Surr)	107		%			70 - 120
Toluene-d8 (Surr)	94		%			80 - 120
4-Bromofluorobenzene (Surr)	92		%			75 - 120

Ms. Liz Scaggs
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Job Number: 560-1385-1

Client Sample ID: R2-D
Lab Sample ID: 560-1385-4

Date Sampled: 08/02/2006 1241
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1613			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U ug/L	0.50	10	1.0
Bis(2-chloroethyl)ether	0.71	U ug/L	0.71	10	1.0
2-Chlorophenol	0.50	U ug/L	0.50	10	1.0
1,3-Dichlorobenzene	0.53	U ug/L	0.53	10	1.0
1,4-Dichlorobenzene	0.74	U ug/L	0.74	10	1.0
Benzyl alcohol	1.4	U ug/L	1.4	20	1.0
1,2-Dichlorobenzene	0.50	U ug/L	0.50	10	1.0
2-Methylphenol	0.50	U ug/L	0.50	10	1.0
Bis(2-chloroisopropyl) ether	0.57	U ug/L	0.57	10	1.0
3 & 4 Methylphenol	0.88	U ug/L	0.88	10	1.0
N-Nitrosodi-n-propylamine	0.65	U ug/L	0.65	10	1.0
Hexachloroethane	0.58	U ug/L	0.58	10	1.0
Nitrobenzene	0.50	U ug/L	0.50	10	1.0
2-Nitrophenol	0.50	U ug/L	0.50	10	1.0
2,4-Dimethylphenol	0.56	U ug/L	0.56	10	1.0
Bis(2-chloroethoxy)methane	0.59	U ug/L	0.59	10	1.0
2,4-Dichlorophenol	0.50	U ug/L	0.50	10	1.0
1,2,4-Trichlorobenzene	0.59	U ug/L	0.59	10	1.0
Naphthalene	0.50	U ug/L	0.50	10	1.0
4-Chloroaniline	0.50	U ug/L	0.50	10	1.0
Hexachlorobutadiene	0.50	U ug/L	0.50	10	1.0
4-Chloro-3-methylphenol	0.50	U ug/L	0.50	10	1.0
2-Methylnaphthalene	0.50	U ug/L	0.50	10	1.0
Hexachlorocyclopentadiene	50	U ug/L	50	50	1.0
2,4,6-Trichlorophenol	0.50	U ug/L	0.50	10	1.0
2,4,5-Trichlorophenol	0.50	U ug/L	0.50	10	1.0
2-Chloronaphthalene	0.50	U ug/L	0.50	10	1.0
2-Nitroaniline	0.50	U ug/L	0.50	50	1.0
Dimethyl phthalate	0.55	U ug/L	0.55	10	1.0
Acenaphthylene	0.50	U ug/L	0.50	10	1.0
2,6-Dinitrotoluene	0.52	U ug/L	0.52	10	1.0
3-Nitroaniline	1.8	U ug/L	1.8	50	1.0
Acenaphthene	0.57	U ug/L	0.57	10	1.0
2,4-Dinitrophenol	20	U ug/L	20	50	1.0
4-Nitrophenol	10	U ug/L	10	50	1.0
2,4-Dinitrotoluene	5.0	U ug/L	5.0	10	1.0
Diethyl phthalate	0.52	U ug/L	0.52	10	1.0
Fluorene	0.61	U ug/L	0.61	10	1.0
4-Chlorophenyl phenyl ether	5.0	U ug/L	5.0	10	1.0
4-Nitroaniline	1.5	U ug/L	1.5	50	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-D
 Lab Sample ID: 560-1385-4

Date Sampled: 08/02/2006 1241
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed: 08/09/2006 1613					
Prep Method: 3520C	Date Prepared: 08/08/2006 1230					
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	U	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0
Surrogate						
2-Fluorophenol	59		%	Acceptance Limits		
Phenol-d5	63		%	10 - 120		
Nitrobenzene-d5	71		%	12 - 120		
2-Fluorobiphenyl	68		%	30 - 120		
2,4,6-Tribromophenol	67		%	26 - 120		
Terphenyl-d14	37		%	25 - 120		
10 - 120						
Method: 8081A	Date Analyzed: 08/14/2006 2112					
Prep Method: 3520C	Date Prepared: 08/08/2006 1700					
alpha-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	U	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	U	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	U	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: R2-D
Lab Sample ID: 560-1385-4

Date Sampled: 08/02/2006 1241
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2112			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
Heptachlor epoxide	0.0028	U ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U ug/L	0.0083	0.050	1.0
Endrin	0.0025	U ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U ug/L	0.050	0.50	1.0
Toxaphene	0.50	U ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U ug/L	0.0027	0.050	1.0
Surrogate				Acceptance Limits	
Tetrachloro-m-xylene	59	%		57 - 127	
DCB Decachlorobiphenyl	22	%		10 - 152	
Method: 8082	Date Analyzed:	08/15/2006 2104			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
Aroclor 1016	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U ug/L	0.17	0.50	1.0
Surrogate				Acceptance Limits	
Tetrachloro-m-xylene	79	%		25 - 140	
DCB Decachlorobiphenyl	32	X %		42 - 133	
Method: 6020	Date Analyzed:	08/09/2006 0229			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	4.6	J ug/L	1.0	5.0	10
Ba	190	B ug/L	1.0	50	10
Cd	1.0	U ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R2-D
Lab Sample ID: 560-1385-4

Date Sampled: 08/02/2006 1241
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0229			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	1.1	J	ug/L	1.1	20
Ni	2.3	J B	ug/L	1.0	5.0
Pb	1.0	U	ug/L	1.0	5.0
Se	1.0	U *	ug/L	1.0	5.0
Zn	53	J	ug/L	50	100
Method: DISS-6020	Date Analyzed:	08/09/2006 0540			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U	ug/L	1.0	5.0
As	4.1	J *	ug/L	1.0	5.0
Ba	180	* B	ug/L	1.0	50
Cd	1.0	U *	ug/L	1.0	5.0
Cr	1.1	U *	ug/L	1.1	20
Ni	2.1	J * B	ug/L	1.0	5.0
Pb	1.0	U	ug/L	1.0	5.0
Se	1.0	U *	ug/L	1.0	5.0
Zn	50	U *	ug/L	50	100
Method: 7470A	Date Analyzed:	08/09/2006 1704			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	U	mg/L	0.00013	0.0020
Method: DISS-7470A	Date Analyzed:	08/09/2006 1724			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	U	mg/L	0.00013	0.0020

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Job Number: 560-1385-1

Client Sample ID: R1-A
 Lab Sample ID: 560-1385-5

Date Sampled: 08/02/2006 1621
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:	08/09/2006 1822			
Prep Method: 5030B	Date Prepared:	08/09/2006 1822			
Chloromethane	0.39	U ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U ug/L	0.20	5.0	1.0
Bromomethane	0.39	U ug/L	0.39	5.0	1.0
Chloroethane	0.40	U ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U ug/L	0.53	50	1.0
Acetone	3.0	J ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U ug/L	0.20	5.0	1.0
Chloroform	0.20	U ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Benzene	0.20	U ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Toluene	0.38	J ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U ug/L	0.20	5.0	1.0
Bromoform	0.50	U ug/L	0.50	5.0	1.0
Styrene	0.50	U ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U ug/L	0.47	5.0	1.0
Xylenes, Total	0.52	J ug/L	0.35	15	1.0
Surrogate				Acceptance Limits	
Dibromofluoromethane (Surr)	102	%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	106	%		70 - 120	
Toluene-d8 (Surr)	97	%		80 - 120	
4-Bromofluorobenzene (Surr)	90	%		75 - 120	

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Job Number: 560-1385-1

Client Sample ID: R1-A
Lab Sample ID: 560-1385-5

Date Sampled: 08/02/2006 1621
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution	
Method: 8270C	Date Analyzed:	08/09/2006 1642				
Prep Method: 3520C	Date Prepared:	08/08/2006 1230				
Phenol	0.50	U	ug/L	0.50	10	1.0
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10	1.0
2-Chlorophenol	0.50	U	ug/L	0.50	10	1.0
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10	1.0
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10	1.0
Benzyl alcohol	1.4	U	ug/L	1.4	20	1.0
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10	1.0
2-Methylphenol	0.50	U	ug/L	0.50	10	1.0
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10	1.0
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10	1.0
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10	1.0
Hexachloroethane	0.58	U	ug/L	0.58	10	1.0
Nitrobenzene	0.50	U	ug/L	0.50	10	1.0
2-Nitrophenol	0.50	U	ug/L	0.50	10	1.0
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10	1.0
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10	1.0
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10	1.0
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10	1.0
Naphthalene	0.50	U	ug/L	0.50	10	1.0
4-Chloroaniline	0.50	U	ug/L	0.50	10	1.0
Hexachlorobutadiene	0.50	U	ug/L	0.50	10	1.0
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10	1.0
2-Methylnaphthalene	0.50	U	ug/L	0.50	10	1.0
Hexachlorocyclopentadiene	50	U	ug/L	50	50	1.0
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10	1.0
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10	1.0
2-Chloronaphthalene	0.50	U	ug/L	0.50	10	1.0
2-Nitroaniline	0.50	U	ug/L	0.50	50	1.0
Dimethyl phthalate	0.55	U	ug/L	0.55	10	1.0
Acenaphthylene	0.50	U	ug/L	0.50	10	1.0
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10	1.0
3-Nitroaniline	1.8	U	ug/L	1.8	50	1.0
Acenaphthene	0.57	U	ug/L	0.57	10	1.0
2,4-Dinitrophenol	20	U	ug/L	20	50	1.0
4-Nitrophenol	10	U	ug/L	10	50	1.0
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10	1.0
Diethyl phthalate	0.52	U	ug/L	0.52	10	1.0
Fluorene	0.61	U	ug/L	0.61	10	1.0
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10	1.0
4-Nitroaniline	1.5	U	ug/L	1.5	50	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-A
 Lab Sample ID: 560-1385-5

Date Sampled: 08/02/2006 1621
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:		08/09/2006 1642			
Prep Method: 3520C	Date Prepared:		08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	U	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0
Surrogate				Acceptance Limits		
2-Fluorophenol	60		%	10 - 120		
Phenol-d5	64		%	12 - 120		
Nitrobenzene-d5	69		%	30 - 120		
2-Fluorobiphenyl	66		%	26 - 120		
2,4,6-Tribromophenol	64		%	25 - 120		
Terphenyl-d14	38		%	10 - 120		

Method: 8081A	Date Analyzed:	08/14/2006 2136
Prep Method: 3520C	Date Prepared:	08/08/2006 1700
alpha-BHC	0.0056	U
beta-BHC	0.0056	U
delta-BHC	0.0025	U
Heptachlor	0.0059	U
Aldrin	0.0025	U

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Job Number: 560-1385-1

Client Sample ID: R1-A
Lab Sample ID: 560-1385-5

Date Sampled: 08/02/2006 1621
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2136				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	59		%	57 - 127		
DCB Decachlorobiphenyl	20		%	10 - 152		
Method: 8082	Date Analyzed:	08/15/2006 2121				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	81		%	25 - 140		
DCB Decachlorobiphenyl	31	X	%	42 - 133		
Method: 6020	Date Analyzed:	08/09/2006 0235				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	4.7	J	ug/L	1.0	5.0	10
Ba	180	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R1-A
Lab Sample ID: 560-1385-5

Date Sampled: 08/02/2006 1621
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed: 08/09/2006 0235					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Cr	1.1	U	ug/L	1.1	20	10
Ni	2.0	J B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U	ug/L	50	100	10
Method: DISS-6020	Date Analyzed: 08/09/2006 0546					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Ag	1.0	U	ug/L	1.0	5.0	10
As	4.8	J *	ug/L	1.0	5.0	10
Ba	210	* B	ug/L	1.0	50	10
Cd	1.0	U *	ug/L	1.0	5.0	10
Cr	1.1	U *	ug/L	1.1	20	10
Ni	2.3	J * B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U *	ug/L	50	100	10
Method: 7470A	Date Analyzed: 08/09/2006 1708					
Prep Method: 7470A	Date Prepared: 08/09/2006 1130					
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed: 08/09/2006 1726					
Prep Method: 7470A	Date Prepared: 08/09/2006 1130					
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-B
 Lab Sample ID: 560-1385-6

Date Sampled: 08/02/2006 1643
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:		08/09/2006 1847			
Prep Method: 5030B	Date Prepared:		08/09/2006 1847			
Chloromethane	0.39	U	ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	0.85	J	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.47	J	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.63	J	ug/L	0.35	15	1.0
Surrogate						Acceptance Limits
Dibromofluoromethane (Surr)	101		%			80 - 120
1,2-Dichloroethane-d4 (Surr)	106		%			70 - 120
Toluene-d8 (Surr)	96		%			80 - 120
4-Bromofluorobenzene (Surr)	91		%			75 - 120

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Job Number: 560-1385-1

Client Sample ID: R1-B
 Lab Sample ID: 560-1385-6

Date Sampled: 08/02/2006 1643
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1710			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	50
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Choronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

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Job Number: 560-1385-1

Client Sample ID: R1-B
 Lab Sample ID: 560-1385-6

Date Sampled: 08/02/2006 1643
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1710			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	ug/L	0.65	10	1.0
Phenanthrene	0.50	ug/L	0.50	10	1.0
Anthracene	0.50	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	ug/L	0.50	10	1.0
Fluoranthene	0.50	ug/L	0.50	10	1.0
Pyrene	0.50	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	ug/L	0.50	10	1.0
Chrysene	0.50	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	ug/L	1.3	10	1.0
Benzoic acid	20	ug/L	20	50	1.0
Surrogate				Acceptance Limits	
2-Fluorophenol	56	%		10 - 120	
Phenol-d5	60	%		12 - 120	
Nitrobenzene-d5	66	%		30 - 120	
2-Fluorobiphenyl	64	%		26 - 120	
2,4,6-Tribromophenol	61	%		25 - 120	
Terphenyl-d14	36	%		10 - 120	
Method: 8081A	Date Analyzed:	08/14/2006 2159			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
alpha-BHC	0.0056	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-B
Lab Sample ID: 560-1385-6

Date Sampled: 08/02/2006 1643
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2159				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	60		%		57 - 127	
DCB Decachlorobiphenyl	18		%		10 - 152	
Method: 8082	Date Analyzed:	08/15/2006 2139				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	80		%		25 - 140	
DCB Decachlorobiphenyl	26	X	%		42 - 133	
Method: 6020	Date Analyzed:	08/09/2006 0241				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	4.5	J	ug/L	1.0	5.0	10
Ba	190	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R1-B
Lab Sample ID: 560-1385-6

Date Sampled: 08/02/2006 1643
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed: 08/09/2006 0241					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Cr	1.1	U	ug/L	1.1	20	10
Ni	1.8	J B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U	ug/L	50	100	10
Method: DISS-6020	Date Analyzed: 08/09/2006 0623					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Ag	1.0	U	ug/L	1.0	5.0	10
As	4.8	J *	ug/L	1.0	5.0	10
Ba	190	* B	ug/L	1.0	50	10
Cd	1.0	U *	ug/L	1.0	5.0	10
Cr	1.1	U *	ug/L	1.1	20	10
Ni	1.2	J * B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U *	ug/L	50	100	10
Method: 7470A	Date Analyzed: 08/11/2006 1554					
Prep Method: 7470A	Date Prepared: 08/11/2006 1105					
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed: 08/11/2006 1623					
Prep Method: 7470A	Date Prepared: 08/11/2006 1105					
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-C
 Lab Sample ID: 560-1385-7

Date Sampled: 08/02/2006 1709
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed: 08/09/2006 1912					
Prep Method: 5030B	Date Prepared: 08/09/2006 1912					
Chloromethane	0.39	U	ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	2.0	J	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.27	J	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.38	J	ug/L	0.35	15	1.0
Surrogate					Acceptance Limits	
Dibromofluoromethane (Surr)	102		%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	108		%		70 - 120	
Toluene-d8 (Surr)	95		%		80 - 120	
4-Bromofluorobenzene (Surr)	92		%		75 - 120	

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Job Number: 560-1385-1

Client Sample ID: R1-C
Lab Sample ID: 560-1385-7

Date Sampled: 08/02/2006 1709
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1738			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	ug/L	0.50	10	1.0
Bis(2-chloroethyl)ether	0.71	ug/L	0.71	10	1.0
2-Chlorophenol	0.50	ug/L	0.50	10	1.0
1,3-Dichlorobenzene	0.53	ug/L	0.53	10	1.0
1,4-Dichlorobenzene	0.74	ug/L	0.74	10	1.0
Benzyl alcohol	1.4	ug/L	1.4	20	1.0
1,2-Dichlorobenzene	0.50	ug/L	0.50	10	1.0
2-Methylphenol	0.50	ug/L	0.50	10	1.0
Bis(2-chloroisopropyl) ether	0.57	ug/L	0.57	10	1.0
3 & 4 Methylphenol	0.88	ug/L	0.88	10	1.0
N-Nitrosodi-n-propylamine	0.65	ug/L	0.65	10	1.0
Hexachloroethane	0.58	ug/L	0.58	10	1.0
Nitrobenzene	0.50	ug/L	0.50	10	1.0
2-Nitrophenol	0.50	ug/L	0.50	10	1.0
2,4-Dimethylphenol	0.56	ug/L	0.56	10	1.0
Bis(2-chloroethoxy)methane	0.59	ug/L	0.59	10	1.0
2,4-Dichlorophenol	0.50	ug/L	0.50	10	1.0
1,2,4-Trichlorobenzene	0.59	ug/L	0.59	10	1.0
Naphthalene	0.50	ug/L	0.50	10	1.0
4-Chloroaniline	0.50	ug/L	0.50	10	1.0
Hexachlorobutadiene	0.50	ug/L	0.50	10	1.0
4-Chloro-3-methylphenol	0.50	ug/L	0.50	10	1.0
2-Methylnaphthalene	0.50	ug/L	0.50	10	1.0
Hexachlorocyclopentadiene	50	ug/L	50	50	1.0
2,4,6-Trichlorophenol	0.50	ug/L	0.50	10	1.0
2,4,5-Trichlorophenol	0.50	ug/L	0.50	10	1.0
2-Chloronaphthalene	0.50	ug/L	0.50	10	1.0
2-Nitroaniline	0.50	ug/L	0.50	50	1.0
Dimethyl phthalate	0.55	ug/L	0.55	10	1.0
Acenaphthylene	0.50	ug/L	0.50	10	1.0
2,6-Dinitrotoluene	0.52	ug/L	0.52	10	1.0
3-Nitroaniline	1.8	ug/L	1.8	50	1.0
Acenaphthene	0.57	ug/L	0.57	10	1.0
2,4-Dinitrophenol	20	ug/L	20	50	1.0
4-Nitrophenol	10	ug/L	10	50	1.0
2,4-Dinitrotoluene	5.0	ug/L	5.0	10	1.0
Diethyl phthalate	0.52	ug/L	0.52	10	1.0
Fluorene	0.61	ug/L	0.61	10	1.0
4-Chlorophenyl phenyl ether	5.0	ug/L	5.0	10	1.0
4-Nitroaniline	1.5	ug/L	1.5	50	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-C
 Lab Sample ID: 560-1385-7

Date Sampled: 08/02/2006 1709
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1738			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	ug/L	0.65	10	1.0
Phenanthrene	0.50	ug/L	0.50	10	1.0
Anthracene	0.50	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	ug/L	0.50	10	1.0
Fluoranthene	0.50	ug/L	0.50	10	1.0
Pyrene	0.50	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	ug/L	0.50	10	1.0
Chrysene	0.50	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	ug/L	1.3	10	1.0
Benzoic acid	20	ug/L	20	50	1.0
Surrogate				Acceptance Limits	
2-Fluorophenol	56	%		10 - 120	
Phenol-d5	60	%		12 - 120	
Nitrobenzene-d5	66	%		30 - 120	
2-Fluorobiphenyl	62	%		26 - 120	
2,4,6-Tribromophenol	59	%		25 - 120	
Terphenyl-d14	37	%		10 - 120	

Method: 8081A	Date Analyzed:	08/14/2006 2223
Prep Method: 3520C	Date Prepared:	08/08/2006 1700
alpha-BHC	0.0056	ug/L
beta-BHC	0.0056	ug/L
delta-BHC	0.0025	ug/L
Heptachlor	0.0059	ug/L
Aldrin	0.0025	ug/L

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Job Number: 560-1385-1

Client Sample ID: R1-C
 Lab Sample ID: 560-1385-7

Date Sampled: 08/02/2006 1709
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2223				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	58		%	57 - 127		
DCB Decachlorobiphenyl	18		%	10 - 152		
Method: 8082	Date Analyzed:	08/15/2006 2156				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	81		%	25 - 140		
DCB Decachlorobiphenyl	27	X	%	42 - 133		
Method: 6020	Date Analyzed:	08/09/2006 0247				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	4.1	J	ug/L	1.0	5.0	10
Ba	170	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R1-C
Lab Sample ID: 560-1385-7

Date Sampled: 08/02/2006 1709
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0247			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	1.1	U ug/L	1.1	20	10
Ni	1.7	J B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0629			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	4.0	J * ug/L	1.0	5.0	10
Ba	180	* B ug/L	1.0	50	10
Cd	1.0	U * ug/L	1.0	5.0	10
Cr	1.1	U * ug/L	1.1	20	10
Ni	1.3	J * B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U * ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/09/2006 1710			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/09/2006 1728			
Prep Method: 7470A	Date Prepared:	08/09/2006 1130			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-D
 Lab Sample ID: 560-1385-8

Date Sampled: 08/02/2006 1729
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:	08/09/2006 1936			
Prep Method: 5030B	Date Prepared:	08/09/2006 1936			
Chloromethane	0.39	U ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U ug/L	0.20	5.0	1.0
Bromomethane	0.39	U ug/L	0.39	5.0	1.0
Chloroethane	0.40	U ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U ug/L	0.53	50	1.0
Acetone	3.4	J ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U ug/L	0.20	5.0	1.0
Chloroform	0.20	U ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Benzene	0.20	U ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Toluene	0.23	J ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U ug/L	0.20	5.0	1.0
Bromoform	0.50	U ug/L	0.50	5.0	1.0
Styrene	0.50	U ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U ug/L	0.47	5.0	1.0
Xylenes, Total	0.36	J ug/L	0.35	15	1.0
Surrogate					
Dibromofluoromethane (Surr)	102	%		Acceptance Limits	
1,2-Dichloroethane-d4 (Surr)	107	%		80 - 120	
Toluene-d8 (Surr)	97	%		70 - 120	
4-Bromofluorobenzene (Surr)	90	%		80 - 120	
				75 - 120	

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Job Number: 560-1385-1

Client Sample ID: R1-D
Lab Sample ID: 560-1385-8

Date Sampled: 08/02/2006 1729
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1806			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U ug/L	0.50	10	1.0
Bis(2-chloroethyl)ether	0.71	U ug/L	0.71	10	1.0
2-Chlorophenol	0.50	U ug/L	0.50	10	1.0
1,3-Dichlorobenzene	0.53	U ug/L	0.53	10	1.0
1,4-Dichlorobenzene	0.74	U ug/L	0.74	10	1.0
Benzyl alcohol	1.4	U ug/L	1.4	20	1.0
1,2-Dichlorobenzene	0.50	U ug/L	0.50	10	1.0
2-Methylphenol	0.50	U ug/L	0.50	10	1.0
Bis(2-chloroisopropyl) ether	0.57	U ug/L	0.57	10	1.0
3 & 4 Methylphenol	0.88	U ug/L	0.88	10	1.0
N-Nitrosodi-n-propylamine	0.65	U ug/L	0.65	10	1.0
Hexachloroethane	0.58	U ug/L	0.58	10	1.0
Nitrobenzene	0.50	U ug/L	0.50	10	1.0
2-Nitrophenol	0.50	U ug/L	0.50	10	1.0
2,4-Dimethylphenol	0.56	U ug/L	0.56	10	1.0
Bis(2-chloroethoxy)methane	0.59	U ug/L	0.59	10	1.0
2,4-Dichlorophenol	0.50	U ug/L	0.50	10	1.0
1,2,4-Trichlorobenzene	0.59	U ug/L	0.59	10	1.0
Naphthalene	0.50	U ug/L	0.50	10	1.0
4-Chloroaniline	0.50	U ug/L	0.50	10	1.0
Hexachlorobutadiene	0.50	U ug/L	0.50	10	1.0
4-Chloro-3-methylphenol	0.50	U ug/L	0.50	10	1.0
2-Methylnaphthalene	0.50	U ug/L	0.50	10	1.0
Hexachlorocyclopentadiene	50	U ug/L	50	50	1.0
2,4,6-Trichlorophenol	0.50	U ug/L	0.50	10	1.0
2,4,5-Trichlorophenol	0.50	U ug/L	0.50	10	1.0
2-Chloronaphthalene	0.50	U ug/L	0.50	10	1.0
2-Nitroaniline	0.50	U ug/L	0.50	50	1.0
Dimethyl phthalate	0.55	U ug/L	0.55	10	1.0
Acenaphthylene	0.50	U ug/L	0.50	10	1.0
2,6-Dinitrotoluene	0.52	U ug/L	0.52	10	1.0
3-Nitroaniline	1.8	U ug/L	1.8	50	1.0
Acenaphthene	0.57	U ug/L	0.57	10	1.0
2,4-Dinitrophenol	20	U ug/L	20	50	1.0
4-Nitrophenol	10	U ug/L	10	50	1.0
2,4-Dinitrotoluene	5.0	U ug/L	5.0	10	1.0
Diethyl phthalate	0.52	U ug/L	0.52	10	1.0
Fluorene	0.61	U ug/L	0.61	10	1.0
4-Chlorophenyl phenyl ether	5.0	U ug/L	5.0	10	1.0
4-Nitroaniline	1.5	U ug/L	1.5	50	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-D
 Lab Sample ID: 560-1385-8

Date Sampled: 08/02/2006 1729
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed: 08/09/2006 1806					
Prep Method: 3520C	Date Prepared: 08/08/2006 1230					
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	U	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0
Surrogate				Acceptance Limits		
2-Fluorophenol	57		%	10 - 120		
Phenol-d5	60		%	12 - 120		
Nitrobenzene-d5	67		%	30 - 120		
2-Fluorobiphenyl	63		%	26 - 120		
2,4,6-Tribromophenol	63		%	25 - 120		
Terphenyl-d14	32		%	10 - 120		
Method: 8081A	Date Analyzed: 08/14/2006 2247					
Prep Method: 3520C	Date Prepared: 08/08/2006 1700					
alpha-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	U	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	U	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	U	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: R1-D
 Lab Sample ID: 560-1385-8

Date Sampled: 08/02/2006 1729
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2247				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	56	X	%		57 - 127	
DCB Decachlorobiphenyl	17		%		10 - 152	
Method: 8082	Date Analyzed:	08/15/2006 2213				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	79		%		25 - 140	
DCB Decachlorobiphenyl	25	X	%		42 - 133	
Method: 6020	Date Analyzed:	08/09/2006 0253				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	5.4		ug/L	1.0	5.0	10
Ba	220	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: R1-D
Lab Sample ID: 560-1385-8

Date Sampled: 08/02/2006 1729
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0253			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	1.1	U	ug/L	1.1	20
Ni	2.3	J B	ug/L	1.0	5.0
Pb	1.0	U	ug/L	1.0	5.0
Se	1.0	U *	ug/L	1.0	5.0
Zn	50	U	ug/L	50	100
Method: DISS-6020	Date Analyzed:	08/09/2006 0635			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U	ug/L	1.0	5.0
As	4.3	J *	ug/L	1.0	5.0
Ba	190	* B	ug/L	1.0	50
Cd	1.0	U *	ug/L	1.0	5.0
Cr	1.1	U *	ug/L	1.1	20
Ni	1.5	J * B	ug/L	1.0	5.0
Pb	1.0	U	ug/L	1.0	5.0
Se	1.0	U *	ug/L	1.0	5.0
Zn	50	U *	ug/L	50	100
Method: 7470A	Date Analyzed:	08/11/2006 1557			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00013	U	mg/L	0.00013	0.0020
Method: DISS-7470A	Date Analyzed:	08/11/2006 1624			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00017	J	mg/L	0.00013	0.0020

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Job Number: 560-1385-1

Client Sample ID: MW-6
 Lab Sample ID: 560-1385-9

Date Sampled: 08/03/2006 0759
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:		08/09/2006 2001			
Prep Method: 5030B	Date Prepared:		08/09/2006 2001			
Chloromethane	0.39	U	ug/L	0.39	5.0	1.0
Vinyl chloride	0.73	J	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	0.46	U	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U*	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.20	U	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U	ug/L	0.35	15	1.0
Surrogate					Acceptance Limits	
Dibromofluoromethane (Surr)	104		%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	107		%		70 - 120	
Toluene-d8 (Surr)	95		%		80 - 120	
4-Bromofluorobenzene (Surr)	92		%		75 - 120	

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Job Number: 560-1385-1

Client Sample ID: MW-6
Lab Sample ID: 560-1385-9

Date Sampled: 08/03/2006 0759
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1834			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	50
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Chloronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

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Job Number: 560-1385-1

Client Sample ID: MW-6
Lab Sample ID: 560-1385-9

Date Sampled: 08/03/2006 0759
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1834			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	ug/L	0.65	10	1.0
Phenanthrene	0.50	ug/L	0.50	10	1.0
Anthracene	0.50	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	ug/L	0.50	10	1.0
Fluoranthene	0.50	ug/L	0.50	10	1.0
Pyrene	0.50	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	ug/L	0.50	10	1.0
Chrysene	0.50	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	ug/L	1.3	10	1.0
Benzoic acid	20	ug/L	20	50	1.0
Surrogate					
2-Fluorophenol	62	%		10 - 120	
Phenol-d5	65	%		12 - 120	
Nitrobenzene-d5	72	%		30 - 120	
2-Fluorobiphenyl	71	%		26 - 120	
2,4,6-Tribromophenol	77	%		25 - 120	
Terphenyl-d14	46	%		10 - 120	
Method: 8081A	Date Analyzed:	08/14/2006 2310			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
alpha-BHC	0.0056	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-6
 Lab Sample ID: 560-1385-9

Date Sampled: 08/03/2006 0759
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2310				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	67		%	57 - 127		
DCB Decachlorobiphenyl	25		%	10 - 152		
Method: 8082	Date Analyzed:	08/15/2006 2230				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	90		%	25 - 140		
DCB Decachlorobiphenyl	36	X	%	42 - 133		
Method: 6020	Date Analyzed:	08/09/2006 0330				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	1.4	J	ug/L	1.0	5.0	10
Ba	1200	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: MW-6
Lab Sample ID: 560-1385-9

Date Sampled: 08/03/2006 0759
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0330			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	3.0	J ug/L	1.1	20	10
Ni	2.1	J B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0641			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	1.0	U * ug/L	1.0	5.0	10
Ba	1000	* B ug/L	1.0	50	10
Cd	1.0	U * ug/L	1.0	5.0	10
Cr	1.1	U * ug/L	1.1	20	10
Ni	1.0	U * ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U * ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/11/2006 1559			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00016	J mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/11/2006 1627			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-31
 Lab Sample ID: 560-1385-10

Date Sampled: 08/03/2006 0920
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:		08/09/2006 2026			
Prep Method: 5030B	Date Prepared:		08/09/2006 2026			
Chloromethane	0.48	J	ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	0.46	U	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U *	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.20	U	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U	ug/L	0.35	15	1.0
Surrogate					Acceptance Limits	
Dibromofluoromethane (Surr)	105		%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	109		%		70 - 120	
Toluene-d8 (Surr)	94		%		80 - 120	
4-Bromofluorobenzene (Surr)	90		%		75 - 120	

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Job Number: 560-1385-1

Client Sample ID: MW-31
Lab Sample ID: 560-1385-10

Date Sampled: 08/03/2006 0920
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006	1902		
Prep Method: 3520C	Date Prepared:	08/08/2006	1230		
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	50
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Chloronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

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Job Number: 560-1385-1

Client Sample ID: MW-31
Lab Sample ID: 560-1385-10

Date Sampled: 08/03/2006 0920
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte		Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C		Date Analyzed:	08/09/2006 1902			
Prep Method: 3520C		Date Prepared:	08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	2.0	J	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0
Surrogate					Acceptance Limits	
2-Fluorophenol	63	%			10 - 120	
Phenol-d5	67	%			12 - 120	
Nitrobenzene-d5	72	%			30 - 120	
2-Fluorobiphenyl	71	%			26 - 120	
2,4,6-Tribromophenol	83	%			25 - 120	
Terphenyl-d14	58	%			10 - 120	

Method: 8081A	Date Analyzed:	08/14/2006 2334
Prep Method: 3520C	Date Prepared:	08/08/2006 1700
alpha-BHC	0.0056	U
beta-BHC	0.0056	U
delta-BHC	0.0025	U
Heptachlor	0.0059	U
Aldrin	0.0025	U

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Job Number: 560-1385-1

Client Sample ID: MW-31
 Lab Sample ID: 560-1385-10

Date Sampled: 08/03/2006 0920
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/14/2006 2334			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
Heptachlor epoxide	0.0028	U ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U ug/L	0.0083	0.050	1.0
Endrin	0.0025	U ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U ug/L	0.050	0.50	1.0
Toxaphene	0.50	U ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U ug/L	0.0027	0.050	1.0
Surrogate					
Tetrachloro-m-xylene	61	%		57 - 127	
DCB Decachlorobiphenyl	15	%		10 - 152	
Method: 8082	Date Analyzed:	08/15/2006 2247			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
Aroclor 1016	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U ug/L	0.17	0.50	1.0
Surrogate					
Tetrachloro-m-xylene	93	%		25 - 140	
DCB Decachlorobiphenyl	19	X %		42 - 133	
Method: 6020	Date Analyzed:	08/09/2006 0337			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	23	ug/L	1.0	5.0	10
Ba	400	B ug/L	1.0	50	10
Cd	1.0	U ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: MW-31
Lab Sample ID: 560-1385-10

Date Sampled: 08/03/2006 0920
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0337			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	1.1	U ug/L	1.1	20	10
Ni	1.0	U ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0648			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	11	U ug/L	1.0	5.0	10
Ba	410	* B ug/L	1.0	50	10
Cd	1.0	U * ug/L	1.0	5.0	10
Cr	1.1	U * ug/L	1.1	20	10
Ni	1.0	U * ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U * ug/L	1.0	5.0	10
Zn	50	U * ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/11/2006 1537			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00023	J mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/11/2006 1601			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0

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 Entact Environmental Services, LLC
 3129 Bass Pro Drive
 Grapevine, TX 76051

Job Number: 560-1385-1

Client Sample ID: MW-34
 Lab Sample ID: 560-1385-11

Date Sampled: 08/03/2006 1040
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:	08/09/2006 2050			
Prep Method: 5030B	Date Prepared:	08/09/2006 2050			
Chloromethane	1.1	J ug/L	0.39	5.0	1.0
Vinyl chloride	1.7	J ug/L	0.20	5.0	1.0
Bromomethane	0.39	U ug/L	0.39	5.0	1.0
Chloroethane	0.40	U ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.91	J ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U ug/L	0.53	50	1.0
Acetone	22	J ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	12	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.24	J ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U * ug/L	0.20	5.0	1.0
Chloroform	0.20	U ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Benzene	67	ug/L	0.20	5.0	1.0
Trichloroethene	0.44	J ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Toluene	0.23	J ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U ug/L	0.50	5.0	1.0
Chlorobenzene	0.52	J ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U ug/L	0.20	5.0	1.0
Bromoform	0.50	U ug/L	0.50	5.0	1.0
Styrene	0.50	U ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	2.0	J ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U ug/L	0.35	15	1.0
Surrogate				Acceptance Limits	
Dibromofluoromethane (Surr)	103	%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	107	%		70 - 120	
Toluene-d8 (Surr)	96	%		80 - 120	
4-Bromofluorobenzene (Surr)	89	%		75 - 120	

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 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1930			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	50
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Chloronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

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 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1930			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	ug/L	0.65	10	1.0
Phenanthrene	0.50	ug/L	0.50	10	1.0
Anthracene	0.50	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	ug/L	0.50	10	1.0
Fluoranthene	0.50	ug/L	0.50	10	1.0
Pyrene	0.50	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	ug/L	0.50	10	1.0
Chrysene	0.50	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	ug/L	1.3	10	1.0
Benzoic acid	20	ug/L	20	50	1.0
Surrogate					
2-Fluorophenol	64	%		10 - 120	
Phenol-d5	68	%		12 - 120	
Nitrobenzene-d5	71	%		30 - 120	
2-Fluorobiphenyl	68	%		26 - 120	
2,4,6-Tribromophenol	83	%		25 - 120	
Terphenyl-d14	47	%		10 - 120	
Acceptance Limits					
Method: 8081A	Date Analyzed:	08/15/2006 0954			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
alpha-BHC	0.0056	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	ug/L	0.0025	0.050	1.0

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 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed: 08/15/2006 0954					
Prep Method: 3520C	Date Prepared: 08/08/2006 1700					
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	67		%		57 - 127	
DCB Decachlorobiphenyl	24		%		10 - 152	
Method: 8082	Date Analyzed: 08/16/2006 1647					
Prep Method: 3520C	Date Prepared: 08/08/2006 1700					
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	98		%		25 - 140	
DCB Decachlorobiphenyl	10	X	%		42 - 133	
Method: 6020	Date Analyzed: 08/09/2006 0343					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Ag	1.0	U	ug/L	1.0	5.0	10
As	5.8		ug/L	1.0	5.0	10
Ba	980	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Date Sampled: 08/03/2006 1040
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0343				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Cr	1.1	U	ug/L	1.1	20	10
Ni	1.0	U	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U	ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0654				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	3.3	J *	ug/L	1.0	5.0	10
Ba	1100	* B	ug/L	1.0	50	10
Cd	1.0	U *	ug/L	1.0	5.0	10
Cr	1.1	U *	ug/L	1.1	20	10
Ni	1.0	U *	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U *	ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/11/2006 1545				
Prep Method: 7470A	Date Prepared:	08/11/2006 1105				
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/11/2006 1602				
Prep Method: 7470A	Date Prepared:	08/11/2006 1105				
Hg	0.00013	J	mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-35
Lab Sample ID: 560-1385-12

Date Sampled: 08/03/2006 1155
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:	08/09/2006 2115			
Prep Method: 5030B	Date Prepared:	08/09/2006 2115			
Chloromethane	0.39	U ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U ug/L	0.20	5.0	1.0
Bromomethane	0.39	U ug/L	0.39	5.0	1.0
Chloroethane	0.40	U ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U ug/L	0.53	50	1.0
Acetone	2.5	J ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U ug/L	0.20	5.0	1.0
Chloroform	0.20	U ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Benzene	0.33	J ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Toluene	0.20	U ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U ug/L	0.20	5.0	1.0
Bromoform	0.50	U ug/L	0.50	5.0	1.0
Styrene	0.50	U ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U ug/L	0.35	15	1.0
Surrogate				Acceptance Limits	
Dibromofluoromethane (Surr)	105	%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	109	%		70 - 120	
Toluene-d8 (Surr)	95	%		80 - 120	
4-Bromofluorobenzene (Surr)	89	%		75 - 120	

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Job Number: 560-1385-1

Client Sample ID: MW-35
 Lab Sample ID: 560-1385-12

Date Sampled: 08/03/2006 1155
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1958			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	50
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Choronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

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Job Number: 560-1385-1

Client Sample ID: MW-35
 Lab Sample ID: 560-1385-12

Date Sampled: 08/03/2006 1155
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 1958			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
4,6-Dinitro-2-methylphenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	ug/L	0.65	10	1.0
Phenanthrene	0.50	ug/L	0.50	10	1.0
Anthracene	0.50	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	ug/L	0.50	10	1.0
Fluoranthene	0.50	ug/L	0.50	10	1.0
Pyrene	0.50	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	ug/L	0.50	10	1.0
Chrysene	0.50	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	7.4	J	1.9	10	1.0
Di-n-octyl phthalate	5.0	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	ug/L	1.3	10	1.0
Benzoic acid	20	U	20	50	1.0
Surrogate					
2-Fluorophenol	61	%		Acceptance Limits	10 - 120
Phenol-d5	65	%			12 - 120
Nitrobenzene-d5	72	%			30 - 120
2-Fluorobiphenyl	68	%			26 - 120
2,4,6-Tribromophenol	79	%			25 - 120
Terphenyl-d14	28	%			10 - 120
Method: 8081A	Date Analyzed:	08/15/2006 1018			
Prep Method: 3520C	Date Prepared:	08/08/2006 1700			
alpha-BHC	0.0056	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	ug/L	0.0025	0.050	1.0
Heptachlor	0.0059	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-35
 Lab Sample ID: 560-1385-12

Date Sampled: 08/03/2006 1155
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/15/2006 1018				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	78		%		57 - 127	
DCB Decachlorobiphenyl	47		%		10 - 152	
Method: 8082	Date Analyzed:	08/16/2006 1704				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	107		%		25 - 140	
DCB Decachlorobiphenyl	57		%		42 - 133	
Method: 6020	Date Analyzed:	08/09/2006 0349				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	1.0	U	ug/L	1.0	5.0	10
Ba	150	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: MW-35
Lab Sample ID: 560-1385-12

Date Sampled: 08/03/2006 1155
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed: 08/09/2006 0349					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Cr	1.1	U	ug/L	1.1	20	10
Ni	2.1	J B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U	ug/L	50	100	10
Method: DISS-6020	Date Analyzed: 08/09/2006 0700					
Prep Method: 3010A	Date Prepared: 08/07/2006 1103					
Ag	1.0	U	ug/L	1.0	5.0	10
As	1.0	U *	ug/L	1.0	5.0	10
Ba	160	* B	ug/L	1.0	50	10
Cd	1.0	U *	ug/L	1.0	5.0	10
Cr	1.1	U *	ug/L	1.1	20	10
Ni	1.0	U *	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	1.0	U *	ug/L	1.0	5.0	10
Zn	50	U *	ug/L	50	100	10
Method: 7470A	Date Analyzed: 08/11/2006 1548					
Prep Method: 7470A	Date Prepared: 08/11/2006 1105					
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed: 08/11/2006 1616					
Prep Method: 7470A	Date Prepared: 08/11/2006 1105					
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-37
 Lab Sample ID: 560-1385-13

Date Sampled: 08/03/2006 1253
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B						
Prep Method: 5030B						
Chloromethane	0.44	J	ug/L	0.39	5.0	1.0
Vinyl chloride	11		ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.22	J	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	4.7	J	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	4.6	J	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U*	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	1.3	J	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	J	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.47	J	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	1.5	J	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.91	J	ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U	ug/L	0.35	15	1.0
Surrogate						Acceptance Limits
Dibromofluoromethane (Surr)	103		%			80 - 120
1,2-Dichloroethane-d4 (Surr)	107		%			70 - 120
Toluene-d8 (Surr)	95		%			80 - 120
4-Bromofluorobenzene (Surr)	91		%			75 - 120

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Job Number: 560-1385-1

Client Sample ID: MW-37
Lab Sample ID: 560-1385-13

Date Sampled: 08/03/2006 1253
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 2026			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U ug/L	0.50	10	1.0
Bis(2-chloroethyl)ether	0.71	U ug/L	0.71	10	1.0
2-Chlorophenol	0.50	U ug/L	0.50	10	1.0
1,3-Dichlorobenzene	0.53	U ug/L	0.53	10	1.0
1,4-Dichlorobenzene	0.74	U ug/L	0.74	10	1.0
Benzyl alcohol	1.4	U ug/L	1.4	20	1.0
1,2-Dichlorobenzene	0.50	U ug/L	0.50	10	1.0
2-Methylphenol	0.50	U ug/L	0.50	10	1.0
Bis(2-chloroisopropyl) ether	0.57	U ug/L	0.57	10	1.0
3 & 4 Methylphenol	0.88	U ug/L	0.88	10	1.0
N-Nitrosodi-n-propylamine	0.65	U ug/L	0.65	10	1.0
Hexachloroethane	0.58	U ug/L	0.58	10	1.0
Nitrobenzene	0.50	U ug/L	0.50	10	1.0
2-Nitrophenol	0.50	U ug/L	0.50	10	1.0
2,4-Dimethylphenol	0.56	U ug/L	0.56	10	1.0
Bis(2-chloroethoxy)methane	0.59	U ug/L	0.59	10	1.0
2,4-Dichlorophenol	0.50	U ug/L	0.50	10	1.0
1,2,4-Trichlorobenzene	0.59	U ug/L	0.59	10	1.0
Naphthalene	0.50	U ug/L	0.50	10	1.0
4-Chloroaniline	0.50	U ug/L	0.50	10	1.0
Hexachlorobutadiene	0.50	U ug/L	0.50	10	1.0
4-Chloro-3-methylphenol	0.50	U ug/L	0.50	10	1.0
2-Methylnaphthalene	0.50	U ug/L	0.50	10	1.0
Hexachlorocyclopentadiene	50	U ug/L	50	50	1.0
2,4,6-Trichlorophenol	0.50	U ug/L	0.50	10	1.0
2,4,5-Trichlorophenol	0.50	U ug/L	0.50	10	1.0
2-Chloronaphthalene	0.50	U ug/L	0.50	10	1.0
2-Nitroaniline	0.50	U ug/L	0.50	50	1.0
Dimethyl phthalate	0.55	U ug/L	0.55	10	1.0
Acenaphthylene	0.50	U ug/L	0.50	10	1.0
2,6-Dinitrotoluene	0.52	U ug/L	0.52	10	1.0
3-Nitroaniline	1.8	U ug/L	1.8	50	1.0
Acenaphthene	0.57	U ug/L	0.57	10	1.0
2,4-Dinitrophenol	20	U ug/L	20	50	1.0
4-Nitrophenol	10	U ug/L	10	50	1.0
2,4-Dinitrotoluene	5.0	U ug/L	5.0	10	1.0
Diethyl phthalate	0.52	U ug/L	0.52	10	1.0
Fluorene	0.61	U ug/L	0.61	10	1.0
4-Chlorophenyl phenyl ether	5.0	U ug/L	5.0	10	1.0
4-Nitroaniline	1.5	U ug/L	1.5	50	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-37
 Lab Sample ID: 560-1385-13

Date Sampled: 08/03/2006 1253
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8270C						
Prep Method: 3520C						
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	1.9	U	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0
Surrogate					Acceptance Limits	
2-Fluorophenol	61		%		10 - 120	
Phenol-d5	64		%		12 - 120	
Nitrobenzene-d5	69		%		30 - 120	
2-Fluorobiphenyl	66		%		26 - 120	
2,4,6-Tribromophenol	82		%		25 - 120	
Terphenyl-d14	44		%		10 - 120	

Method: 8081A	Date Analyzed:	08/15/2006	1041
Prep Method: 3520C	Date Prepared:	08/08/2006	1700
alpha-BHC	0.0056	U	ug/L
beta-BHC	0.0056	U	ug/L
delta-BHC	0.0025	U	ug/L
Heptachlor	0.0059	U	ug/L
Aldrin	0.0025	U	ug/L

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Job Number: 560-1385-1

Client Sample ID: MW-37
Lab Sample ID: 560-1385-13

Date Sampled: 08/03/2006 1253
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/15/2006 1041				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.0089	U	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	62		%		57 - 127	
DCB Decachlorobiphenyl	14		%		10 - 152	
Method: 8082	Date Analyzed:	08/16/2006 1722				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate					Acceptance Limits	
Tetrachloro-m-xylene	94		%		25 - 140	
DCB Decachlorobiphenyl	16	X	%		42 - 133	
Method: 6020	Date Analyzed:	08/09/2006 0355				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	4.0	J	ug/L	1.0	5.0	10
Ba	490	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: MW-37
Lab Sample ID: 560-1385-13

Date Sampled: 08/03/2006 1253
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:		08/09/2006 0355			
Prep Method: 3010A	Date Prepared:		08/07/2006 1103			
Cr	1.1	U	ug/L	1.1	20	10
Ni	1.4	J B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	6.2	*	ug/L	1.0	5.0	10
Zn	50	U	ug/L	50	100	10
Method: DISS-6020	Date Analyzed:		08/09/2006 0706			
Prep Method: 3010A	Date Prepared:		08/07/2006 1103			
Ag	1.0	U	ug/L	1.0	5.0	10
As	3.4	J *	ug/L	1.0	5.0	10
Ba	450	* B	ug/L	1.0	50	10
Cd	1.0	U *	ug/L	1.0	5.0	10
Cr	1.1	U *	ug/L	1.1	20	10
Ni	2.5	J * B	ug/L	1.0	5.0	10
Pb	1.0	U	ug/L	1.0	5.0	10
Se	5.3	*	ug/L	1.0	5.0	10
Zn	50	U *	ug/L	50	100	10
Method: 7470A	Date Analyzed:		08/11/2006 1550			
Prep Method: 7470A	Date Prepared:		08/11/2006 1105			
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:		08/11/2006 1619			
Prep Method: 7470A	Date Prepared:		08/11/2006 1105			
Hg	0.00013	U	mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-39
 Lab Sample ID: 560-1385-14

Date Sampled: 08/03/2006 1358
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8260B			Date Analyzed:	08/09/2006 2205		
Prep Method: 5030B			Date Prepared:	08/09/2006 2205		
Chloromethane	0.51	J	ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U	ug/L	0.20	5.0	1.0
Bromomethane	0.39	U	ug/L	0.39	5.0	1.0
Chloroethane	0.40	U	ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U	ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U	ug/L	0.53	50	1.0
Acetone	4.8	J	ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U	ug/L	0.20	5.0	1.0
Chloroform	0.20	U	ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U	ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Benzene	0.20	U	ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U	ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U	ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U	ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Toluene	0.20	U	ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U	ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U	ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U	ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U	ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U	ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U	ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U	ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U	ug/L	0.20	5.0	1.0
Bromoform	0.50	U	ug/L	0.50	5.0	1.0
Styrene	0.50	U	ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U	ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U	ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U	ug/L	0.35	15	1.0
Surrogate					Acceptance Limits	
Dibromofluoromethane (Surr)	103		%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	109		%		70 - 120	
Toluene-d8 (Surr)	96		%		80 - 120	
4-Bromofluorobenzene (Surr)	90		%		75 - 120	

Ms. Liz Scaggs
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3129 Bass Pro Drive
Grapevine, TX 76051

Job Number: 560-1385-1

Client Sample ID: MW-39
Lab Sample ID: 560-1385-14

Date Sampled: 08/03/2006 1358
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8270C	Date Analyzed:	08/09/2006 2055			
Prep Method: 3520C	Date Prepared:	08/08/2006 1230			
Phenol	0.50	U	ug/L	0.50	10
Bis(2-chloroethyl)ether	0.71	U	ug/L	0.71	10
2-Chlorophenol	0.50	U	ug/L	0.50	10
1,3-Dichlorobenzene	0.53	U	ug/L	0.53	10
1,4-Dichlorobenzene	0.74	U	ug/L	0.74	10
Benzyl alcohol	1.4	U	ug/L	1.4	20
1,2-Dichlorobenzene	0.50	U	ug/L	0.50	10
2-Methylphenol	0.50	U	ug/L	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	ug/L	0.57	10
3 & 4 Methylphenol	0.88	U	ug/L	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	ug/L	0.65	10
Hexachloroethane	0.58	U	ug/L	0.58	10
Nitrobenzene	0.50	U	ug/L	0.50	10
2-Nitrophenol	0.50	U	ug/L	0.50	10
2,4-Dimethylphenol	0.56	U	ug/L	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	ug/L	0.59	10
2,4-Dichlorophenol	0.50	U	ug/L	0.50	10
1,2,4-Trichlorobenzene	0.59	U	ug/L	0.59	10
Naphthalene	0.50	U	ug/L	0.50	10
4-Chloroaniline	0.50	U	ug/L	0.50	10
Hexachlorobutadiene	0.50	U	ug/L	0.50	10
4-Chloro-3-methylphenol	0.50	U	ug/L	0.50	10
2-Methylnaphthalene	0.50	U	ug/L	0.50	10
Hexachlorocyclopentadiene	50	U	ug/L	50	50
2,4,6-Trichlorophenol	0.50	U	ug/L	0.50	10
2,4,5-Trichlorophenol	0.50	U	ug/L	0.50	10
2-Chloronaphthalene	0.50	U	ug/L	0.50	10
2-Nitroaniline	0.50	U	ug/L	0.50	50
Dimethyl phthalate	0.55	U	ug/L	0.55	10
Acenaphthylene	0.50	U	ug/L	0.50	10
2,6-Dinitrotoluene	0.52	U	ug/L	0.52	10
3-Nitroaniline	1.8	U	ug/L	1.8	50
Acenaphthene	0.57	U	ug/L	0.57	10
2,4-Dinitrophenol	20	U	ug/L	20	50
4-Nitrophenol	10	U	ug/L	10	50
2,4-Dinitrotoluene	5.0	U	ug/L	5.0	10
Diethyl phthalate	0.52	U	ug/L	0.52	10
Fluorene	0.61	U	ug/L	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	ug/L	5.0	10
4-Nitroaniline	1.5	U	ug/L	1.5	50

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Job Number: 560-1385-1

Client Sample ID: MW-39
Lab Sample ID: 560-1385-14

Date Sampled: 08/03/2006 1358
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8270C			Date Analyzed:	08/09/2006 2055		
Prep Method: 3520C			Date Prepared:	08/08/2006 1230		
4,6-Dinitro-2-methylphenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodiphenylamine	0.51	U	ug/L	0.51	10	1.0
4-Bromophenyl phenyl ether	0.74	U	ug/L	0.74	10	1.0
Hexachlorobenzene	0.65	U	ug/L	0.65	10	1.0
Phenanthrene	0.50	U	ug/L	0.50	10	1.0
Anthracene	0.50	U	ug/L	0.50	10	1.0
Di-n-butyl phthalate	0.50	U	ug/L	0.50	10	1.0
Fluoranthene	0.50	U	ug/L	0.50	10	1.0
Pyrene	0.50	U	ug/L	0.50	10	1.0
Butyl benzyl phthalate	0.50	U	ug/L	0.50	10	1.0
Benzo[a]anthracene	0.50	U	ug/L	0.50	10	1.0
Chrysene	0.50	U	ug/L	0.50	10	1.0
Bis(2-ethylhexyl) phthalate	9.6	J	ug/L	1.9	10	1.0
Di-n-octyl phthalate	5.0	U	ug/L	5.0	10	1.0
Benzo[b]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[k]fluoranthene	0.50	U	ug/L	0.50	10	1.0
Benzo[a]pyrene	0.50	U	ug/L	0.50	10	1.0
Indeno[1,2,3-cd]pyrene	0.50	U	ug/L	0.50	10	1.0
Dibenz(a,h)anthracene	0.50	U	ug/L	0.50	10	1.0
Benzo[g,h,i]perylene	0.50	U	ug/L	0.50	10	1.0
3,3'-Dichlorobenzidine	0.50	U	ug/L	0.50	20	1.0
Pentachlorophenol	5.0	U	ug/L	5.0	50	1.0
N-Nitrosodimethylamine	1.3	U	ug/L	1.3	10	1.0
Benzoic acid	20	U	ug/L	20	50	1.0
Surrogate						Acceptance Limits
2-Fluorophenol	61		%			10 - 120
Phenol-d5	66		%			12 - 120
Nitrobenzene-d5	70		%			30 - 120
2-Fluorobiphenyl	68		%			26 - 120
2,4,6-Tribromophenol	81		%			25 - 120
Terphenyl-d14	51		%			10 - 120
Method: 8081A			Date Analyzed:	08/15/2006 1105		
Prep Method: 3520C			Date Prepared:	08/08/2006 1700		
alpha-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
beta-BHC	0.0056	U	ug/L	0.0056	0.050	1.0
delta-BHC	0.0025	U	ug/L	0.0025	0.050	1.0
Heptachlor	0.0081	J	ug/L	0.0059	0.050	1.0
Aldrin	0.0025	U	ug/L	0.0025	0.050	1.0

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Job Number: 560-1385-1

Client Sample ID: MW-39
Lab Sample ID: 560-1385-14

Date Sampled: 08/03/2006 1358
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier		Unit	MDL	RL	Dilution
Method: 8081A	Date Analyzed:	08/15/2006 1105				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Heptachlor epoxide	0.0028	U	ug/L	0.0028	0.050	1.0
4,4'-DDE	0.0026	U	ug/L	0.0026	0.050	1.0
Endosulfan I	0.020	J	ug/L	0.0089	0.050	1.0
Dieldrin	0.0083	U	ug/L	0.0083	0.050	1.0
Endrin	0.0025	U	ug/L	0.0025	0.050	1.0
4,4'-DDD	0.0029	U	ug/L	0.0029	0.050	1.0
Endosulfan II	0.0035	U	ug/L	0.0035	0.050	1.0
4,4'-DDT	0.0034	U	ug/L	0.0034	0.050	1.0
Methoxychlor	0.023	U	ug/L	0.023	0.050	1.0
Endosulfan sulfate	0.0039	U	ug/L	0.0039	0.050	1.0
Endrin ketone	0.0073	U	ug/L	0.0073	0.050	1.0
Chlordane (technical)	0.050	U	ug/L	0.050	0.50	1.0
Toxaphene	0.50	U	ug/L	0.50	5.0	1.0
gamma-BHC (Lindane)	0.0027	U	ug/L	0.0027	0.050	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	56	X	%	57 - 127		
DCB Decachlorobiphenyl	17		%	10 - 152		
Method: 8082	Date Analyzed:	08/16/2006 1739				
Prep Method: 3520C	Date Prepared:	08/08/2006 1700				
Aroclor 1016	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1221	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1232	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1242	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1248	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1254	0.17	U	ug/L	0.17	0.50	1.0
Aroclor 1260	0.17	U	ug/L	0.17	0.50	1.0
Surrogate				Acceptance Limits		
Tetrachloro-m-xylene	100		%	25 - 140		
DCB Decachlorobiphenyl	22	X	%	42 - 133		
Method: 6020	Date Analyzed:	08/09/2006 0401				
Prep Method: 3010A	Date Prepared:	08/07/2006 1103				
Ag	1.0	U	ug/L	1.0	5.0	10
As	51		ug/L	1.0	5.0	10
Ba	240	B	ug/L	1.0	50	10
Cd	1.0	U	ug/L	1.0	5.0	10

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Job Number: 560-1385-1

Client Sample ID: MW-39
Lab Sample ID: 560-1385-14

Date Sampled: 08/03/2006 1358
Date Received: 08/04/2006 1040
Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 6020	Date Analyzed:	08/09/2006 0401			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Cr	1.1	U ug/L	1.1	20	10
Ni	12	B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U ug/L	1.0	5.0	10
Zn	50	U ug/L	50	100	10
Method: DISS-6020	Date Analyzed:	08/09/2006 0712			
Prep Method: 3010A	Date Prepared:	08/07/2006 1103			
Ag	1.0	U ug/L	1.0	5.0	10
As	51	* ug/L	1.0	5.0	10
Ba	320	* B ug/L	1.0	50	10
Cd	1.0	U * ug/L	1.0	5.0	10
Cr	1.1	U * ug/L	1.1	20	10
Ni	1.8	J * B ug/L	1.0	5.0	10
Pb	1.0	U ug/L	1.0	5.0	10
Se	1.0	U ug/L	1.0	5.0	10
Zn	50	U * ug/L	50	100	10
Method: 7470A	Date Analyzed:	08/11/2006 1551			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0
Method: DISS-7470A	Date Analyzed:	08/11/2006 1621			
Prep Method: 7470A	Date Prepared:	08/11/2006 1105			
Hg	0.00013	U mg/L	0.00013	0.0020	1.0

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Job Number: 560-1385-1

Client Sample ID: TRIP BLANK
 Lab Sample ID: 560-1385-15

Date Sampled: 08/03/2006 0000
 Date Received: 08/04/2006 1040
 Client Matrix: Water

Analyte	Result/Qualifier	Unit	MDL	RL	Dilution
Method: 8260B	Date Analyzed:	08/07/2006 1811			
Prep Method: 5030B	Date Prepared:	08/07/2006 1811			
Chloromethane	0.39	U ug/L	0.39	5.0	1.0
Vinyl chloride	0.20	U ug/L	0.20	5.0	1.0
Bromomethane	0.39	U ug/L	0.39	5.0	1.0
Chloroethane	0.40	U ug/L	0.40	5.0	1.0
1,1-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
Carbon disulfide	0.20	U ug/L	0.20	5.0	1.0
Methylene Chloride	0.53	U ug/L	0.53	50	1.0
Acetone	0.46	U ug/L	0.46	100	1.0
trans-1,2-Dichloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1-Dichloroethane	0.20	U ug/L	0.20	5.0	1.0
Vinyl acetate	0.20	U ug/L	0.20	5.0	1.0
Chloroform	0.20	U ug/L	0.20	5.0	1.0
Carbon tetrachloride	0.25	U ug/L	0.25	5.0	1.0
1,1,1-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Benzene	0.20	U ug/L	0.20	5.0	1.0
Trichloroethene	0.32	U ug/L	0.32	5.0	1.0
1,2-Dichloropropane	0.20	U ug/L	0.20	5.0	1.0
Bromodichloromethane	0.20	U ug/L	0.20	5.0	1.0
cis-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Toluene	0.20	U ug/L	0.20	5.0	1.0
methyl isobutyl ketone	0.50	U ug/L	0.50	5.0	1.0
trans-1,3-Dichloropropene	0.20	U ug/L	0.20	5.0	1.0
Tetrachloroethene	0.20	U ug/L	0.20	5.0	1.0
1,1,2-Trichloroethane	0.20	U ug/L	0.20	5.0	1.0
Chlorodibromomethane	0.22	U ug/L	0.22	5.0	1.0
2-Hexanone	0.50	U ug/L	0.50	5.0	1.0
Chlorobenzene	0.20	U ug/L	0.20	5.0	1.0
Ethylbenzene	0.20	U ug/L	0.20	5.0	1.0
Bromoform	0.50	U ug/L	0.50	5.0	1.0
Styrene	0.50	U ug/L	0.50	5.0	1.0
1,1,2,2-Tetrachloroethane	0.20	U ug/L	0.20	5.0	1.0
Methyl Ethyl Ketone	0.47	U ug/L	0.47	5.0	1.0
Xylenes, Total	0.35	U ug/L	0.35	15	1.0
Surrogate				Acceptance Limits	
Dibromofluoromethane (Surr)	106	%		80 - 120	
1,2-Dichloroethane-d4 (Surr)	115	%		70 - 120	
Toluene-d8 (Surr)	101	%		80 - 120	
4-Bromofluorobenzene (Surr)	94	%		75 - 120	

DATA REPORTING QUALIFIERS

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Section	Qualifier	Description
GC/MS VOA	B	Compound was found in the blank and sample.
	U	Indicates the analyte was analyzed for but not detected.
	*	LCS or LCSD exceeds the control limits
	J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.
GC/MS Semi VOA	U	Indicates the analyte was analyzed for but not detected.
	F	MS or MSD exceeds the control limits
	J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.
GC Semi VOA	U	Indicates the analyte was analyzed for but not detected.
	F	MS or MSD exceeds the control limits
	J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.
	X	Surrogate exceeds the control limits
Metals	B	Compound was found in the blank and sample.
	U	Indicates the analyte was analyzed for but not detected.
	*	LCS or LCSD exceeds the control limits
	J	Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value.

QUALITY CONTROL RESULTS

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS VOA					
Analysis Batch:560-3834					
LCS 560-3834/1	Lab Control Spike	T	Water	8260B	
MB 560-3834/2	Method Blank	T	Water	8260B	
560-1385-1	R2-A	T	Water	8260B	
560-1385-2	R2-B	T	Water	8260B	
560-1385-15TB	TRIP BLANK	T	Water	8260B	
Analysis Batch:560-3940					
LCS 560-3940/1	Lab Control Spike	T	Water	8260B	
MB 560-3940/2	Method Blank	T	Water	8260B	
560-1385-3	R2-C	T	Water	8260B	
560-1385-3MS	Matrix Spike	T	Water	8260B	
560-1385-3MSD	Matrix Spike Duplicate	T	Water	8260B	
560-1385-4	R2-D	T	Water	8260B	
560-1385-5	R1-A	T	Water	8260B	
560-1385-6	R1-B	T	Water	8260B	
560-1385-7	R1-C	T	Water	8260B	
560-1385-8	R1-D	T	Water	8260B	
560-1385-9	MW-6	T	Water	8260B	
560-1385-10	MW-31	T	Water	8260B	
560-1385-11	MW-34	T	Water	8260B	
560-1385-12	MW-35	T	Water	8260B	
560-1385-13	MW-37	T	Water	8260B	
560-1385-14	MW-39	T	Water	8260B	

Report Basis

T = Total

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC/MS Semi VOA					
Prep Batch: 560-3816					
LCS 560-3816/2-A	Lab Control Spike	T	Water	3520C	
MB 560-3816/1-A	Method Blank	T	Water	3520C	
560-1385-1	R2-A	T	Water	3520C	
560-1385-1MS	Matrix Spike	T	Water	3520C	
560-1385-1MSD	Matrix Spike Duplicate	T	Water	3520C	
560-1385-2	R2-B	T	Water	3520C	
560-1385-3	R2-C	T	Water	3520C	
560-1385-4	R2-D	T	Water	3520C	
560-1385-5	R1-A	T	Water	3520C	
560-1385-6	R1-B	T	Water	3520C	
560-1385-7	R1-C	T	Water	3520C	
560-1385-8	R1-D	T	Water	3520C	
560-1385-9	MW-6	T	Water	3520C	
560-1385-10	MW-31	T	Water	3520C	
560-1385-11	MW-34	T	Water	3520C	
560-1385-12	MW-35	T	Water	3520C	
560-1385-13	MW-37	T	Water	3520C	
560-1385-14	MW-39	T	Water	3520C	
Analysis Batch: 560-3854					
LCS 560-3816/2-A	Lab Control Spike	T	Water	8270C	560-3816
MB 560-3816/1-A	Method Blank	T	Water	8270C	560-3816
560-1385-1	R2-A	T	Water	8270C	560-3816
560-1385-1MS	Matrix Spike	T	Water	8270C	560-3816
560-1385-1MSD	Matrix Spike Duplicate	T	Water	8270C	560-3816
560-1385-2	R2-B	T	Water	8270C	560-3816
560-1385-3	R2-C	T	Water	8270C	560-3816
560-1385-4	R2-D	T	Water	8270C	560-3816
560-1385-5	R1-A	T	Water	8270C	560-3816
560-1385-6	R1-B	T	Water	8270C	560-3816
560-1385-7	R1-C	T	Water	8270C	560-3816
560-1385-8	R1-D	T	Water	8270C	560-3816
560-1385-9	MW-6	T	Water	8270C	560-3816
560-1385-10	MW-31	T	Water	8270C	560-3816
560-1385-11	MW-34	T	Water	8270C	560-3816
560-1385-12	MW-35	T	Water	8270C	560-3816
560-1385-13	MW-37	T	Water	8270C	560-3816
560-1385-14	MW-39	T	Water	8270C	560-3816

Report Basis

T = Total

STL Corpus Christi

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Prep Batch: 560-3831					
LCS 560-3831/2-A	Lab Control Spike	T	Water	3520C	
MB 560-3831/1-A	Method Blank	T	Water	3520C	
560-1385-1	R2-A	T	Water	3520C	
560-1385-2	R2-B	T	Water	3520C	
560-1385-2MS	Matrix Spike	T	Water	3520C	
560-1385-2MSD	Matrix Spike Duplicate	T	Water	3520C	
560-1385-3	R2-C	T	Water	3520C	
560-1385-4	R2-D	T	Water	3520C	
560-1385-5	R1-A	T	Water	3520C	
560-1385-6	R1-B	T	Water	3520C	
560-1385-7	R1-C	T	Water	3520C	
560-1385-8	R1-D	T	Water	3520C	
560-1385-9	MW-6	T	Water	3520C	
560-1385-10	MW-31	T	Water	3520C	
560-1385-11	MW-34	T	Water	3520C	
560-1385-12	MW-35	T	Water	3520C	
560-1385-13	MW-37	T	Water	3520C	
560-1385-14	MW-39	T	Water	3520C	
Prep Batch: 560-3832					
LCS 560-3832/2-B	Lab Control Spike	T	Water	3520C	
MB 560-3832/1-A	Method Blank	T	Water	3520C	
560-1385-1	R2-A	T	Water	3520C	
560-1385-2	R2-B	T	Water	3520C	
560-1385-3	R2-C	T	Water	3520C	
560-1385-3MS	Matrix Spike	T	Water	3520C	
560-1385-3MSD	Matrix Spike Duplicate	T	Water	3520C	
560-1385-4	R2-D	T	Water	3520C	
560-1385-5	R1-A	T	Water	3520C	
560-1385-6	R1-B	T	Water	3520C	
560-1385-7	R1-C	T	Water	3520C	
560-1385-8	R1-D	T	Water	3520C	
560-1385-9	MW-6	T	Water	3520C	
560-1385-10	MW-31	T	Water	3520C	
560-1385-11	MW-34	T	Water	3520C	
560-1385-12	MW-35	T	Water	3520C	
560-1385-13	MW-37	T	Water	3520C	
560-1385-14	MW-39	T	Water	3520C	

Quality Control Results

Client: Enact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
GC Semi VOA					
Analysis Batch:560-3996					
LCS 560-3832/2-B	Lab Control Spike	T	Water	8081A	560-3832
MB 560-3832/1-A	Method Blank	T	Water	8081A	560-3832
560-1385-1	R2-A	T	Water	8081A	560-3832
560-1385-2	R2-B	T	Water	8081A	560-3832
560-1385-3	R2-C	T	Water	8081A	560-3832
560-1385-3MS	Matrix Spike	T	Water	8081A	560-3832
560-1385-3MSD	Matrix Spike Duplicate	T	Water	8081A	560-3832
560-1385-4	R2-D	T	Water	8081A	560-3832
560-1385-5	R1-A	T	Water	8081A	560-3832
560-1385-6	R1-B	T	Water	8081A	560-3832
560-1385-7	R1-C	T	Water	8081A	560-3832
560-1385-8	R1-D	T	Water	8081A	560-3832
560-1385-9	MW-6	T	Water	8081A	560-3832
560-1385-10	MW-31	T	Water	8081A	560-3832
560-1385-11	MW-34	T	Water	8081A	560-3832
560-1385-12	MW-35	T	Water	8081A	560-3832
560-1385-13	MW-37	T	Water	8081A	560-3832
560-1385-14	MW-39	T	Water	8081A	560-3832
Analysis Batch:560-4015					
LCS 560-3831/2-A	Lab Control Spike	T	Water	8082	560-3831
MB 560-3831/1-A	Method Blank	T	Water	8082	560-3831
560-1385-1	R2-A	T	Water	8082	560-3831
560-1385-2	R2-B	T	Water	8082	560-3831
560-1385-2MS	Matrix Spike	T	Water	8082	560-3831
560-1385-2MSD	Matrix Spike Duplicate	T	Water	8082	560-3831
560-1385-3	R2-C	T	Water	8082	560-3831
560-1385-4	R2-D	T	Water	8082	560-3831
560-1385-5	R1-A	T	Water	8082	560-3831
560-1385-6	R1-B	T	Water	8082	560-3831
560-1385-7	R1-C	T	Water	8082	560-3831
560-1385-8	R1-D	T	Water	8082	560-3831
560-1385-9	MW-6	T	Water	8082	560-3831
560-1385-10	MW-31	T	Water	8082	560-3831
560-1385-11	MW-34	T	Water	8082	560-3831
560-1385-12	MW-35	T	Water	8082	560-3831
560-1385-13	MW-37	T	Water	8082	560-3831
560-1385-14	MW-39	T	Water	8082	560-3831

Report Basis

T = Total

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
Metals					
Prep Batch: 560-3757					
LCS 560-3757/26-A	Lab Control Spike	D	Water	3010A	
LCS 560-3757/2-A	Lab Control Spike	T	Water	3010A	
MB 560-3757/25-A	Method Blank	D	Water	3010A	
MB 560-3757/1-A	Method Blank	T	Water	3010A	
560-1385-1	R2-A	D	Water	3010A	
560-1385-1	R2-A	T	Water	3010A	
560-1385-1MS	Matrix Spike	D	Water	3010A	
560-1385-1MSD	Matrix Spike Duplicate	D	Water	3010A	
560-1385-2	R2-B	D	Water	3010A	
560-1385-2	R2-B	T	Water	3010A	
560-1385-3	R2-C	D	Water	3010A	
560-1385-3	R2-C	T	Water	3010A	
560-1385-4	R2-D	D	Water	3010A	
560-1385-4	R2-D	T	Water	3010A	
560-1385-5	R1-A	D	Water	3010A	
560-1385-5	R1-A	T	Water	3010A	
560-1385-6	R1-B	D	Water	3010A	
560-1385-6	R1-B	T	Water	3010A	
560-1385-7	R1-C	D	Water	3010A	
560-1385-7	R1-C	T	Water	3010A	
560-1385-8	R1-D	D	Water	3010A	
560-1385-8	R1-D	T	Water	3010A	
560-1385-9	MW-6	D	Water	3010A	
560-1385-9	MW-6	T	Water	3010A	
560-1385-10	MW-31	D	Water	3010A	
560-1385-10	MW-31	T	Water	3010A	
560-1385-11	MW-34	D	Water	3010A	
560-1385-11	MW-34	T	Water	3010A	
560-1385-12	MW-35	D	Water	3010A	
560-1385-12	MW-35	T	Water	3010A	
560-1385-13	MW-37	D	Water	3010A	
560-1385-13	MW-37	T	Water	3010A	
560-1385-14	MW-39	D	Water	3010A	
560-1385-14	MW-39	T	Water	3010A	

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
Metals					
Prep Batch: 560-3833					
LCS 560-3833/18-A	Lab Control Spike	D	Water	7470A	
MB 560-3833/19-A	Method Blank	D	Water	7470A	
560-1385-1	R2-A	D	Water	7470A	
560-1385-1	R2-A	T	Water	7470A	
560-1385-2	R2-B	D	Water	7470A	
560-1385-2	R2-B	T	Water	7470A	
560-1385-3	R2-C	D	Water	7470A	
560-1385-3	R2-C	T	Water	7470A	
560-1385-4	R2-D	D	Water	7470A	
560-1385-4	R2-D	T	Water	7470A	
560-1385-5	R1-A	D	Water	7470A	
560-1385-5	R1-A	T	Water	7470A	
560-1385-7	R1-C	D	Water	7470A	
560-1385-7	R1-C	T	Water	7470A	
Analysis Batch: 560-3844					
LCS 560-3833/18-A	Lab Control Spike	D	Water	7470A	560-3833
MB 560-3833/19-A	Method Blank	D	Water	7470A	560-3833
560-1385-1	R2-A	D	Water	7470A	560-3833
560-1385-1	R2-A	T	Water	7470A	560-3833
560-1385-2	R2-B	D	Water	7470A	560-3833
560-1385-2	R2-B	T	Water	7470A	560-3833
560-1385-3	R2-C	D	Water	7470A	560-3833
560-1385-3	R2-C	T	Water	7470A	560-3833
560-1385-4	R2-D	D	Water	7470A	560-3833
560-1385-4	R2-D	T	Water	7470A	560-3833
560-1385-5	R1-A	D	Water	7470A	560-3833
560-1385-5	R1-A	T	Water	7470A	560-3833
560-1385-7	R1-C	D	Water	7470A	560-3833
560-1385-7	R1-C	T	Water	7470A	560-3833

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
Metals					
Analysis Batch:560-3885					
LCS 560-3757/26-A	Lab Control Spike	D	Water	6020	560-3757
LCS 560-3757/2-A	Lab Control Spike	T	Water	6020	560-3757
MB 560-3757/25-A	Method Blank	D	Water	6020	560-3757
MB 560-3757/1-A	Method Blank	T	Water	6020	560-3757
560-1385-1	R2-A	D	Water	6020	560-3757
560-1385-1	R2-A	T	Water	6020	560-3757
560-1385-1MS	Matrix Spike	D	Water	6020	560-3757
560-1385-1MSD	Matrix Spike Duplicate	D	Water	6020	560-3757
560-1385-2	R2-B	D	Water	6020	560-3757
560-1385-2	R2-B	T	Water	6020	560-3757
560-1385-3	R2-C	D	Water	6020	560-3757
560-1385-3	R2-C	T	Water	6020	560-3757
560-1385-4	R2-D	D	Water	6020	560-3757
560-1385-4	R2-D	T	Water	6020	560-3757
560-1385-5	R1-A	D	Water	6020	560-3757
560-1385-5	R1-A	T	Water	6020	560-3757
560-1385-6	R1-B	D	Water	6020	560-3757
560-1385-6	R1-B	T	Water	6020	560-3757
560-1385-7	R1-C	D	Water	6020	560-3757
560-1385-7	R1-C	T	Water	6020	560-3757
560-1385-8	R1-D	D	Water	6020	560-3757
560-1385-8	R1-D	T	Water	6020	560-3757
560-1385-9	MW-6	D	Water	6020	560-3757
560-1385-9	MW-6	T	Water	6020	560-3757
560-1385-10	MW-31	D	Water	6020	560-3757
560-1385-10	MW-31	T	Water	6020	560-3757
560-1385-11	MW-34	D	Water	6020	560-3757
560-1385-11	MW-34	T	Water	6020	560-3757
560-1385-12	MW-35	D	Water	6020	560-3757
560-1385-12	MW-35	T	Water	6020	560-3757
560-1385-13	MW-37	D	Water	6020	560-3757
560-1385-13	MW-37	T	Water	6020	560-3757
560-1385-14	MW-39	D	Water	6020	560-3757
560-1385-14	MW-39	T	Water	6020	560-3757

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
Metals					
Prep Batch: 560-3912					
LCS 560-3912/3-A	Lab Control Spike	T	Water	7470A	
MB 560-3912/4-A	Method Blank	D	Water	7470A	
560-1385-6	R1-B	D	Water	7470A	
560-1385-6	R1-B	T	Water	7470A	
560-1385-8	R1-D	D	Water	7470A	
560-1385-8	R1-D	T	Water	7470A	
560-1385-9	MW-6	D	Water	7470A	
560-1385-9	MW-6	T	Water	7470A	
560-1385-10	MW-31	D	Water	7470A	
560-1385-10	MW-31	T	Water	7470A	
560-1385-10MS	Matrix Spike	D	Water	7470A	
560-1385-10MSD	Matrix Spike Duplicate	T	Water	7470A	
560-1385-11	MW-34	D	Water	7470A	
560-1385-11	MW-34	T	Water	7470A	
560-1385-12	MW-35	D	Water	7470A	
560-1385-12	MW-35	T	Water	7470A	
560-1385-13	MW-37	D	Water	7470A	
560-1385-13	MW-37	T	Water	7470A	
560-1385-14	MW-39	D	Water	7470A	
560-1385-14	MW-39	T	Water	7470A	
Analysis Batch: 560-3921					
LCS 560-3912/3-A	Lab Control Spike	T	Water	7470A	560-3912
MB 560-3912/4-A	Method Blank	D	Water	7470A	560-3912
560-1385-6	R1-B	D	Water	7470A	560-3912
560-1385-6	R1-B	T	Water	7470A	560-3912
560-1385-8	R1-D	D	Water	7470A	560-3912
560-1385-8	R1-D	T	Water	7470A	560-3912
560-1385-9	MW-6	D	Water	7470A	560-3912
560-1385-9	MW-6	T	Water	7470A	560-3912
560-1385-10	MW-31	D	Water	7470A	560-3912
560-1385-10	MW-31	T	Water	7470A	560-3912
560-1385-10MS	Matrix Spike	D	Water	7470A	560-3912
560-1385-10MSD	Matrix Spike Duplicate	T	Water	7470A	560-3912
560-1385-11	MW-34	D	Water	7470A	560-3912
560-1385-11	MW-34	T	Water	7470A	560-3912
560-1385-12	MW-35	D	Water	7470A	560-3912
560-1385-12	MW-35	T	Water	7470A	560-3912
560-1385-13	MW-37	D	Water	7470A	560-3912
560-1385-13	MW-37	T	Water	7470A	560-3912
560-1385-14	MW-39	D	Water	7470A	560-3912
560-1385-14	MW-39	T	Water	7470A	560-3912

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

QC Association Summary

Lab Sample ID	Client Sample ID	Report Basis	Client Matrix	Method	Prep Batch
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Report Basis

D = Dissolved

T = Total

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Surrogate Recovery Report

8260B Volatile Organic Compounds by GC/MS

Client Matrix: Water

Lab Sample ID	Client Sample	(12DCE) (%Rec)	(BFB) (%Rec)	(DBFM) (%Rec)	(TOL) (%Rec)
560-1385-1	R2-A	113	92	109	99
560-1385-2	R2-B	113	94	108	101
560-1385-3	R2-C	104	92	101	95
560-1385-4	R2-D	107	92	103	94
560-1385-5	R1-A	106	90	102	97
560-1385-6	R1-B	106	91	101	96
560-1385-7	R1-C	108	92	102	95
560-1385-8	R1-D	107	90	102	97
560-1385-9	MW-6	107	92	104	95
560-1385-10	MW-31	109	90	105	94
560-1385-11	MW-34	107	89	103	96
560-1385-12	MW-35	109	89	105	95
560-1385-13	MW-37	107	91	103	95
560-1385-14	MW-39	109	90	103	96
560-1385-15TB	TRIP BLANK	115	94	106	101
560-1385-3MS	R2-C	103	89	103	94
560-1385-3MSD	R2-C	106	91	101	94
LCS 560-3834/1		100	98	106	100
LCS 560-3940/1		100	93	102	96
MB 560-3834/2		105	97	103	100
MB 560-3940/2		102	92	98	97

Surrogate	Acceptance Limits
(12DCE)	1,2-Dichloroethane-d4 (Surr)
(BFB)	4-Bromofluorobenzene (Surr)
(DBFM)	Dibromofluoromethane (Surr)
(TOL)	Toluene-d8 (Surr)

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Surrogate Recovery Report

8270C Semivolatile Compounds by Gas Chromatography/Mass Spectrometry (GC/MS)

Client Matrix: Water

<u>Lab Sample ID</u>	<u>Client Sample</u>	(2FP) (%Rec)	(FBP) (%Rec)	(NBZ) (%Rec)	(PHL) (%Rec)	(TBP) (%Rec)	(TPH) (%Rec)
560-1385-1	R2-A	56	59	66	60	63	49
560-1385-2	R2-B	62	70	73	66	68	40
560-1385-3	R2-C	54	62	63	58	58	34
560-1385-4	R2-D	59	68	71	63	67	37
560-1385-5	R1-A	60	66	69	64	64	38
560-1385-6	R1-B	56	64	66	60	61	36
560-1385-7	R1-C	56	62	66	60	59	37
560-1385-8	R1-D	57	63	67	60	63	32
560-1385-9	MW-6	62	71	72	65	77	46
560-1385-10	MW-31	63	71	72	67	83	58
560-1385-11	MW-34	64	68	71	68	83	47
560-1385-12	MW-35	61	68	72	65	79	28
560-1385-13	MW-37	61	66	69	64	82	44
560-1385-14	MW-39	61	68	70	66	81	51
560-1385-1MS	R2-A	47	50	54	50	57	28
560-1385-1MSD	R2-A	53	55	59	56	70	35
LCS 560-3816/2-A		61	64	68	66	84	85
MB 560-3816/1-A		69	75	79	73	84	86

Surrogate		Acceptance Limits
(2FP)	2-Fluorophenol	10 - 120
(FBP)	2-Fluorobiphenyl	26 - 120
(NBZ)	Nitrobenzene-d5	30 - 120
(PHL)	Phenol-d5	12 - 120
(TBP)	2,4,6-Tribromophenol	25 - 120
(TPH)	Terphenyl-d14	10 - 120

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Surrogate Recovery Report

8081A Organochlorine Pesticides by Gas Chromatography

Client Matrix: Water

Lab Sample ID	Client Sample	(DCB 1) (%Rec)	(TCX 1) (%Rec)
560-1385-1	R2-A	19	60
560-1385-2	R2-B	17	58
560-1385-3	R2-C	19	60
560-1385-4	R2-D	22	59
560-1385-5	R1-A	20	59
560-1385-6	R1-B	18	60
560-1385-7	R1-C	18	58
560-1385-8	R1-D	17	56 X
560-1385-9	MW-6	25	67
560-1385-10	MW-31	15	61
560-1385-11	MW-34	24	67
560-1385-12	MW-35	47	78
560-1385-13	MW-37	14	62
560-1385-14	MW-39	17	56 X
560-1385-3MS	R2-C	28	63
560-1385-3MSD	R2-C	24	60
LCS 560-3832/2-B		66	77
MB 560-3832/1-A		63	73

Surrogate

Acceptance Limits

(DCB 1)	DCB Decachlorobiphenyl	10 - 152
(TCX 1)	Tetrachloro-m-xylene	57 - 127

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Surrogate Recovery Report

8082 Polychlorinated Biphenyls (PCBs) by Gas Chromatography

Client Matrix: Water

Lab Sample ID	Client Sample	(DCB 1) (%Rec)	(TCX 1) (%Rec)
560-1385-1	R2-A	27 X	86
560-1385-2	R2-B	26 X	83
560-1385-3	R2-C	27 X	82
560-1385-4	R2-D	32 X	79
560-1385-5	R1-A	31 X	81
560-1385-6	R1-B	26 X	80
560-1385-7	R1-C	27 X	81
560-1385-8	R1-D	25 X	79
560-1385-9	MW-6	36 X	90
560-1385-10	MW-31	19 X	93
560-1385-11	MW-34	10 X	98
560-1385-12	MW-35	57	107
560-1385-13	MW-37	16 X	94
560-1385-14	MW-39	22 X	100
560-1385-2MS	R2-B	24 X	81
560-1385-2MSD	R2-B	33 X	83
LCS 560-3831/2-A		83	100
MB 560-3831/1-A		82	96

Surrogate

Acceptance Limits

(DCB 1)	DCB Decachlorobiphenyl	42 - 133
(TCX 1)	Tetrachloro-m-xylene	25 - 140

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3834

Lab Sample ID: MB 560-3834/2
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/07/2006 1136
 Date Prepared: 08/07/2006 1136

Analysis Batch: 560-3834
 Prep Batch: N/A
 Units: ug/L

Method: 8260B
Preparation: 5030B

Instrument ID: Hewlett Packard GCMS
 Lab File ID: 08070606.D
 Initial Weight/Volume: 5 mL
 Final Weight/Volume: 5 mL

Analyte	Result	Qual	MDL	RL
Chloromethane	0.39	U	0.39	5.0
Vinyl chloride	0.20	U	0.20	5.0
Bromomethane	0.39	U	0.39	5.0
Chloroethane	0.40	U	0.40	5.0
1,1-Dichloroethene	0.20	U	0.20	5.0
Carbon disulfide	0.20	U	0.20	5.0
Methylene Chloride	0.64	J	0.53	50
Acetone	0.46	U	0.46	100
trans-1,2-Dichloroethene	0.20	U	0.20	5.0
1,1-Dichloroethane	0.20	U	0.20	5.0
Vinyl acetate	0.20	U	0.20	5.0
Chloroform	0.20	U	0.20	5.0
Carbon tetrachloride	0.25	U	0.25	5.0
1,1,1-Trichloroethane	0.20	U	0.20	5.0
Benzene	0.20	U	0.20	5.0
Trichloroethene	0.32	U	0.32	5.0
1,2-Dichloropropane	0.20	U	0.20	5.0
Bromodichloromethane	0.20	U	0.20	5.0
cis-1,3-Dichloropropene	0.20	U	0.20	5.0
Toluene	0.20	U	0.20	5.0
methyl isobutyl ketone	0.50	U	0.50	5.0
trans-1,3-Dichloropropene	0.20	U	0.20	5.0
Tetrachloroethene	0.20	U	0.20	5.0
1,1,2-Trichloroethane	0.20	U	0.20	5.0
Chlorodibromomethane	0.22	U	0.22	5.0
2-Hexanone	0.50	U	0.50	5.0
Chlorobenzene	0.20	U	0.20	5.0
Ethylbenzene	0.20	U	0.20	5.0
Bromoform	0.50	U	0.50	5.0
Styrene	0.50	U	0.50	5.0
1,1,2,2-Tetrachloroethane	0.20	U	0.20	5.0
Methyl Ethyl Ketone	0.47	U	0.47	5.0
Xylenes, Total	0.35	U	0.35	15

Surrogate	% Rec	Acceptance Limits
Dibromofluoromethane (Surr)	103	80 - 120
1,2-Dichloroethane-d4 (Surr)	105	70 - 120
Toluene-d8 (Surr)	100	80 - 120
4-Bromofluorobenzene (Surr)	97	75 - 120

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Control Spike - Batch: 560-3834

Method: 8260B

Preparation: 5030B

Lab Sample ID: LCS 560-3834/1
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/07/2006 1022
 Date Prepared: 08/07/2006 1022

Analysis Batch: 560-3834
 Prep Batch: N/A
 Units: ug/L

Instrument ID: Hewlett Packard GCMS
 Lab File ID: 08070603.D
 Initial Weight/Volume: 5 mL
 Final Weight/Volume: 5 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Chloromethane	50.0	58	116	40 - 125	
Vinyl chloride	50.0	55	110	50 - 145	
Bromomethane	50.0	50	100	30 - 145	
Chloroethane	50.0	54	109	60 - 135	
1,1-Dichloroethene	50.0	54	108	70 - 130	
Carbon disulfide	50.0	56	111	35 - 160	
Methylene Chloride	50.0	57	115	55 - 140	B
Acetone	50.0	52	105	40 - 140	J
trans-1,2-Dichloroethene	50.0	54	108	60 - 140	
1,1-Dichloroethane	50.0	54	109	70 - 135	
Vinyl acetate	50.0	71	142	80 - 148	
Chloroform	50.0	53	107	65 - 135	
Carbon tetrachloride	50.0	52	104	65 - 140	
1,1,1-Trichloroethane	50.0	53	106	65 - 130	
Benzene	50.0	50	100	80 - 120	
Trichloroethene	50.0	44	89	70 - 125	
1,2-Dichloropropane	50.0	53	106	75 - 125	
Bromodichloromethane	50.0	52	104	75 - 120	
cis-1,3-Dichloropropene	50.0	44	87	70 - 130	
Toluene	50.0	49	98	75 - 120	
methyl isobutyl ketone	50.0	51	102	60 - 135	
trans-1,3-Dichloropropene	50.0	56	111	55 - 140	
Tetrachloroethene	50.0	44	89	45 - 150	
1,1,2-Trichloroethane	50.0	51	101	75 - 125	
Chlorodibromomethane	50.0	46	92	60 - 135	
2-Hexanone	50.0	50	99	55 - 130	
Chlorobenzene	50.0	46	92	80 - 120	
Ethylbenzene	50.0	48	96	75 - 125	
Bromoform	50.0	45	91	70 - 130	
Styrene	50.0	50	99	65 - 135	
1,1,2,2-Tetrachloroethane	50.0	52	105	65 - 130	
Methyl Ethyl Ketone	50.0	56	111	30 - 150	
Xylenes, Total	150	140	96	80 - 120	
Surrogate		% Rec		Acceptance Limits	
Dibromofluoromethane (Surr)		106		80 - 120	
1,2-Dichloroethane-d4 (Surr)		100		70 - 120	
Toluene-d8 (Surr)		100		80 - 120	
4-Bromofluorobenzene (Surr)		98		75 - 120	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3940

Lab Sample ID: MB 560-3940/2
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/09/2006 1552
 Date Prepared: 08/09/2006 1552

Analysis Batch: 560-3940
 Prep Batch: N/A
 Units: ug/L

Method: 8260B
Preparation: 5030B

Instrument ID: Hewlett Packard GCMS
 Lab File ID: 08090606.D
 Initial Weight/Volume: 5 mL
 Final Weight/Volume: 5 mL

Analyte	Result	Qual	MDL	RL
Chloromethane	0.39	U	0.39	5.0
Vinyl chloride	0.20	U	0.20	5.0
Bromomethane	0.39	U	0.39	5.0
Chloroethane	0.40	U	0.40	5.0
1,1-Dichloroethene	0.20	U	0.20	5.0
Carbon disulfide	0.20	U	0.20	5.0
Methylene Chloride	1.4	J	0.53	50
Acetone	0.46	U	0.46	100
trans-1,2-Dichloroethene	0.20	U	0.20	5.0
1,1-Dichloroethane	0.20	U	0.20	5.0
Vinyl acetate	0.20	U	0.20	5.0
Chloroform	0.20	U	0.20	5.0
Carbon tetrachloride	0.25	U	0.25	5.0
1,1,1-Trichloroethane	0.20	U	0.20	5.0
Benzene	0.20	U	0.20	5.0
Trichloroethene	0.32	U	0.32	5.0
1,2-Dichloropropane	0.20	U	0.20	5.0
Bromodichloromethane	0.20	U	0.20	5.0
cis-1,3-Dichloropropene	0.20	U	0.20	5.0
Toluene	0.20	U	0.20	5.0
methyl isobutyl ketone	0.50	U	0.50	5.0
trans-1,3-Dichloropropene	0.20	U	0.20	5.0
Tetrachloroethene	0.20	U	0.20	5.0
1,1,2-Trichloroethane	0.20	U	0.20	5.0
Chlorodibromomethane	0.22	U	0.22	5.0
2-Hexanone	0.50	U	0.50	5.0
Chlorobenzene	0.20	U	0.20	5.0
Ethylbenzene	0.20	U	0.20	5.0
Bromoform	0.50	U	0.50	5.0
Styrene	0.50	U	0.50	5.0
1,1,2,2-Tetrachloroethane	0.20	U	0.20	5.0
Methyl Ethyl Ketone	0.47	U	0.47	5.0
Xylenes, Total	0.35	U	0.35	15

Surrogate	% Rec	Acceptance Limits
Dibromofluoromethane (Surr)	98	80 - 120
1,2-Dichloroethane-d4 (Surr)	102	70 - 120
Toluene-d8 (Surr)	97	80 - 120
4-Bromofluorobenzene (Surr)	92	75 - 120

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Control Spike - Batch: 560-3940

Method: 8260B

Preparation: 5030B

Lab Sample ID: LCS 560-3940/1

Client Matrix: Water

Dilution: 1.0

Date Analyzed: 08/09/2006 1438

Date Prepared: 08/09/2006 1438

Analysis Batch: 560-3940

Prep Batch: N/A

Units: ug/L

Instrument ID: Hewlett Packard GCMS

Lab File ID: 08090603.D

Initial Weight/Volume: 5 mL

Final Weight/Volume: 5 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Chloromethane	50.0	59	117	40 - 125	
Vinyl chloride	50.0	54	109	50 - 145	
Bromomethane	50.0	51	101	30 - 145	
Chloroethane	50.0	55	110	60 - 135	
1,1-Dichloroethene	50.0	54	109	70 - 130	
Carbon disulfide	50.0	56	112	35 - 160	
Methylene Chloride	50.0	58	116	55 - 140	B
Acetone	50.0	53	107	40 - 140	J
trans-1,2-Dichloroethene	50.0	56	112	60 - 140	
1,1-Dichloroethane	50.0	56	113	70 - 135	
Vinyl acetate	50.0	75	150	80 - 148	*
Chloroform	50.0	55	110	65 - 135	
Carbon tetrachloride	50.0	53	106	65 - 140	
1,1,1-Trichloroethane	50.0	54	109	65 - 130	
Benzene	50.0	52	104	80 - 120	
Trichloroethene	50.0	46	91	70 - 125	
1,2-Dichloropropane	50.0	54	108	75 - 125	
Bromodichloromethane	50.0	54	109	75 - 120	
cis-1,3-Dichloropropene	50.0	45	91	70 - 130	
Toluene	50.0	50	100	75 - 120	
methyl isobutyl ketone	50.0	55	109	60 - 135	
trans-1,3-Dichloropropene	50.0	58	115	55 - 140	
Tetrachloroethene	50.0	46	92	45 - 150	
1,1,2-Trichloroethane	50.0	52	105	75 - 125	
Chlorodibromomethane	50.0	48	96	60 - 135	
2-Hexanone	50.0	52	105	55 - 130	
Chlorobenzene	50.0	47	93	80 - 120	
Ethylbenzene	50.0	49	98	75 - 125	
Bromoform	50.0	47	93	70 - 130	
Styrene	50.0	50	100	65 - 135	
1,1,2,2-Tetrachloroethane	50.0	56	111	65 - 130	
Methyl Ethyl Ketone	50.0	58	117	30 - 150	
Xylenes, Total	150	150	98	80 - 120	
Surrogate		% Rec		Acceptance Limits	
Dibromofluoromethane (Surr)		102		80 - 120	
1,2-Dichloroethane-d4 (Surr)		100		70 - 120	
Toluene-d8 (Surr)		96		80 - 120	
4-Bromofluorobenzene (Surr)		93		75 - 120	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 560-3940

Method: 8260B
Preparation: 5030B

MS Lab Sample ID: 560-1385-3 Analysis Batch: 560-3940
Client Matrix: Water Prep Batch: N/A
Dilution: 1.0
Date Analyzed: 08/09/2006 1644
Date Prepared: 08/09/2006 1644

Instrument ID: Hewlett Packard GCMS
Lab File ID: 08090608.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

MSD Lab Sample ID: 560-1385-3 Analysis Batch: 560-3940
Client Matrix: Water Prep Batch: N/A
Dilution: 1.0
Date Analyzed: 08/09/2006 1708
Date Prepared: 08/09/2006 1708

Instrument ID: Hewlett Packard GCMS
Lab File ID: 08090609.D
Initial Weight/Volume: 5 mL
Final Weight/Volume: 5 mL

Analyte	% Rec.		RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD				
1,1-Dichloroethene	105	111	70 - 130	5.7	20.0	
Benzene	100	107	80 - 120	6.7	20.0	
Trichloroethene	88	92	70 - 125	4.7	20.0	
Toluene	94	101	75 - 120	7.7	20.0	
Chlorobenzene	85	91	80 - 120	7.1	20.0	
Surrogate	MS % Rec		MSD % Rec		Acceptance Limits	
Dibromofluoromethane (Surr)	103		101		80 - 120	
1,2-Dichloroethane-d4 (Surr)	103		106		70 - 120	
Toluene-d8 (Surr)	94		94		80 - 120	
4-Bromofluorobenzene (Surr)	89		91		75 - 120	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3816

Method: 8270C

Preparation: 3520C

Lab Sample ID: MB 560-3816/1-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1229
Date Prepared: 08/08/2006 1230

Analysis Batch: 560-3854
Prep Batch: 560-3816
Units: ug/L

Instrument ID: Agilent GCMS [Method 827
Lab File ID: 08090605.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

Analyte	Result	Qual	MDL	RL
Phenol	0.50	U	0.50	10
Bis(2-chloroethyl)ether	0.71	U	0.71	10
2-Chlorophenol	0.50	U	0.50	10
1,3-Dichlorobenzene	0.53	U	0.53	10
1,4-Dichlorobenzene	0.74	U	0.74	10
Benzyl alcohol	1.4	U	1.4	20
1,2-Dichlorobenzene	0.50	U	0.50	10
2-Methylphenol	0.50	U	0.50	10
Bis(2-chloroisopropyl) ether	0.57	U	0.57	10
3 & 4 Methylphenol	0.88	U	0.88	10
N-Nitrosodi-n-propylamine	0.65	U	0.65	10
Hexachloroethane	0.58	U	0.58	10
Nitrobenzene	0.50	U	0.50	10
2-Nitrophenol	0.50	U	0.50	10
2,4-Dimethylphenol	0.56	U	0.56	10
Bis(2-chloroethoxy)methane	0.59	U	0.59	10
2,4-Dichlorophenol	0.50	U	0.50	10
1,2,4-Trichlorobenzene	0.59	U	0.59	10
Naphthalene	0.50	U	0.50	10
4-Chloroaniline	0.50	U	0.50	10
Hexachlorobutadiene	0.50	U	0.50	10
4-Chloro-3-methylphenol	0.50	U	0.50	10
2-Methylnaphthalene	0.50	U	0.50	10
Hexachlorocyclopentadiene	50	U	50	50
2,4,6-Trichlorophenol	0.50	U	0.50	10
2,4,5-Trichlorophenol	0.50	U	0.50	10
2-Chloronaphthalene	0.50	U	0.50	10
2-Nitroaniline	0.50	U	0.50	50
Dimethyl phthalate	0.55	U	0.55	10
Acenaphthylene	0.50	U	0.50	10
2,6-Dinitrotoluene	0.52	U	0.52	10
3-Nitroaniline	1.8	U	1.8	50
Acenaphthene	0.57	U	0.57	10
2,4-Dinitrophenol	20	U	20	50
4-Nitrophenol	10	U	10	50
2,4-Dinitrotoluene	5.0	U	5.0	10
Diethyl phthalate	0.52	U	0.52	10
Fluorene	0.61	U	0.61	10
4-Chlorophenyl phenyl ether	5.0	U	5.0	10
4-Nitroaniline	1.5	U	1.5	50
4,6-Dinitro-2-methylphenol	5.0	U	5.0	50

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3816

Lab Sample ID: MB 560-3816/1-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1229
Date Prepared: 08/08/2006 1230

Analysis Batch: 560-3854
Prep Batch: 560-3816
Units: ug/L

Method: 8270C
Preparation: 3520C

Instrument ID: Agilent GCMS [Method 827
Lab File ID: 08090605.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

Analyte	Result	Qual	MDL	RL
N-Nitrosodiphenylamine	0.51	U	0.51	10
4-Bromophenyl phenyl ether	0.74	U	0.74	10
Hexachlorobenzene	0.65	U	0.65	10
Phenanthrene	0.50	U	0.50	10
Anthracene	0.50	U	0.50	10
Di-n-butyl phthalate	0.50	U	0.50	10
Fluoranthene	0.50	U	0.50	10
Pyrene	0.50	U	0.50	10
Butyl benzyl phthalate	0.50	U	0.50	10
Benzo[a]anthracene	0.50	U	0.50	10
Chrysene	0.50	U	0.50	10
Bis(2-ethylhexyl) phthalate	1.9	U	1.9	10
Di-n-octyl phthalate	5.0	U	5.0	10
Benzo[b]fluoranthene	0.50	U	0.50	10
Benzo[k]fluoranthene	0.50	U	0.50	10
Benzo[a]pyrene	0.50	U	0.50	10
Indeno[1,2,3-cd]pyrene	0.50	U	0.50	10
Dibenz(a,h)anthracene	0.50	U	0.50	10
Benzo[g,h,i]perylene	0.50	U	0.50	10
3,3'-Dichlorobenzidine	0.50	U	0.50	20
Pentachlorophenol	5.0	U	5.0	50
N-Nitrosodimethylamine	1.3	U	1.3	10
Benzoic acid	20	U	20	50

Surrogate	% Rec	Acceptance Limits
2-Fluorophenol	69	10 - 120
Phenol-d5	73	12 - 120
Nitrobenzene-d5	79	30 - 120
2-Fluorobiphenyl	75	26 - 120
2,4,6-Tribromophenol	84	25 - 120
Terphenyl-d14	86	10 - 120

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Control Spike - Batch: 560-3816

Method: 8270C

Preparation: 3520C

Lab Sample ID: LCS 560-3816/2-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1257
Date Prepared: 08/08/2006 1230

Analysis Batch: 560-3854
Prep Batch: 560-3816
Units: ug/L

Instrument ID: Agilent GCMS [Method 827
Lab File ID: 08090606.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Phenol	100	68.2	68	20 - 120	
Bis(2-chloroethyl)ether	100	65.1	65	35 - 110	
2-Chlorophenol	100	65.7	66	35 - 105	
1,3-Dichlorobenzene	100	47.1	47	30 - 100	
1,4-Dichlorobenzene	100	49.0	49	30 - 100	
Benzyl alcohol	100	73.3	73	30 - 110	
1,2-Dichlorobenzene	100	50.7	51	35 - 100	
2-Methylphenol	100	70.7	71	40 - 110	
Bis(2-chloroisopropyl) ether	100	64.3	64	25 - 130	
3 & 4 Methylphenol	200	131	66	30 - 110	
N-Nitrosodi-n-propylamine	100	65.6	66	35 - 130	
Hexachloroethane	100	46.5	46	30 - 95	
Nitrobenzene	100	69.7	70	45 - 110	
2-Nitrophenol	100	70.5	71	40 - 115	
2,4-Dimethylphenol	100	68.5	69	30 - 110	
Bis(2-chloroethoxy)methane	100	70.6	71	45 - 105	
2,4-Dichlorophenol	100	73.2	73	50 - 105	
1,2,4-Trichlorobenzene	100	65.9	66	35 - 105	
Naphthalene	100	68.4	68	40 - 100	
4-Chloroaniline	100	57.0	57	15 - 110	
Hexachlorobutadiene	100	62.2	62	25 - 105	
4-Chloro-3-methylphenol	100	72.9	73	45 - 110	
2-Methylnaphthalene	100	69.2	69	45 - 105	
Hexachlorocyclopentadiene	100	43.6	44	10 - 120	U
2,4,6-Trichlorophenol	100	74.0	74	50 - 115	
2,4,5-Trichlorophenol	100	75.8	76	50 - 110	
2-Chloronaphthalene	100	71.6	72	50 - 105	
2-Nitroaniline	100	75.6	76	50 - 115	
Dimethyl phthalate	100	79.7	80	25 - 125	
Acenaphthylene	100	73.7	74	50 - 105	
2,6-Dinitrotoluene	100	77.2	77	50 - 115	
3-Nitroaniline	100	76.4	76	20 - 125	
Acenaphthene	100	75.0	75	45 - 110	
2,4-Dinitrophenol	100	94.2	94	15 - 140	
4-Nitrophenol	100	91.8	92	20 - 120	
2,4-Dinitrotoluene	100	83.4	83	50 - 120	
Diethyl phthalate	100	84.2	84	40 - 120	
Fluorene	100	77.3	77	50 - 110	
4-Chlorophenyl phenyl ether	100	76.4	76	50 - 110	
4-Nitroaniline	100	81.3	81	35 - 120	
4,6-Dinitro-2-methylphenol	100	93.3	93	40 - 130	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Enact Environmental Services, LLC

Job Number: 560-1385-1

Lab Control Spike - Batch: 560-3816

Method: 8270C

Preparation: 3520C

Lab Sample ID: LCS 560-3816/2-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1257
Date Prepared: 08/08/2006 1230

Analysis Batch: 560-3854
Prep Batch: 560-3816
Units: ug/L

Instrument ID: Agilent GCMS [Method 827
Lab File ID: 08090606.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
N-Nitrosodiphenylamine	100	77.4	77	50 - 110	
4-Bromophenyl phenyl ether	100	81.9	82	50 - 115	
Hexachlorobenzene	100	84.8	85	50 - 110	
Phenanthrene	100	85.4	85	50 - 115	
Anthracene	100	86.5	86	55 - 110	
Di-n-butyl phthalate	100	88.0	88	55 - 115	
Fluoranthene	100	86.9	87	55 - 115	
Pyrene	100	88.4	88	50 - 130	
Butyl benzyl phthalate	100	86.7	87	45 - 115	
Benzo[a]anthracene	100	90.2	90	55 - 110	
Chrysene	100	88.0	88	55 - 110	
Bis(2-ethylhexyl) phthalate	100	87.4	87	40 - 125	
Di-n-octyl phthalate	100	91.3	91	35 - 135	
Benzo[b]fluoranthene	100	92.2	92	45 - 120	
Benzo[k]fluoranthene	100	79.8	80	45 - 125	
Benzo[a]pyrene	100	87.2	87	55 - 110	
Indeno[1,2,3-cd]pyrene	100	91.7	92	45 - 125	
Dibenz(a,h)anthracene	100	90.0	90	40 - 125	
Benzo[g,h,i]perylene	100	88.5	89	40 - 125	
3,3'-Dichlorobenzidine	100	83.6	84	20 - 110	
Pentachlorophenol	100	86.3	86	40 - 115	
N-Nitrosodimethylamine	100	59.0	59	25 - 110	
Benzoic acid	100	56.5	56	10 - 121	
Surrogate		% Rec		Acceptance Limits	
2-Fluorophenol		61		10 - 120	
Phenol-d5		66		12 - 120	
Nitrobenzene-d5		68		30 - 120	
2-Fluorobiphenyl		64		26 - 120	
2,4,6-Tribromophenol		84		25 - 120	
Terphenyl-d14		85		10 - 120	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 560-3816

Method: 8270C

Preparation: 3520C

MS Lab Sample ID: 560-1385-1 Analysis Batch: 560-3854
Client Matrix: Water Prep Batch: 560-3816
Dilution: 1.0
Date Analyzed: 08/09/2006 1325
Date Prepared: 08/08/2006 1230

Instrument ID: Agilent GCMS [Method
Lab File ID: 08090607.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

MSD Lab Sample ID: 560-1385-1 Analysis Batch: 560-3854
Client Matrix: Water Prep Batch: 560-3816
Dilution: 1.0
Date Analyzed: 08/09/2006 1353
Date Prepared: 08/08/2006 1230

Instrument ID: Agilent GCMS [Method 827
Lab File ID: 08090608.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

Analyte	% Rec.		RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD				
Phenol	51	58	20 - 120	12	20	
Bis(2-chloroethyl)ether	51	58	35 - 110	13	20	
2-Chlorophenol	51	59	35 - 105	14	20	
1,3-Dichlorobenzene	41	46	30 - 100	13	20	
1,4-Dichlorobenzene	42	48	30 - 100	13	20	
Benzyl alcohol	55	64	30 - 110	14	20	
1,2-Dichlorobenzene	43	49	35 - 100	13	20	
2-Methylphenol	50	59	40 - 110	16	20	
Bis(2-chloroisopropyl) ether	50	55	25 - 130	10	20	
3 & 4 Methylphenol	49	57	30 - 110	15	20	
N-Nitrosodi-n-propylamine	51	57	35 - 130	12	20	
Hexachloroethane	40	45	30 - 95	12	20	
Nitrobenzene	53	61	45 - 110	13	20	
2-Nitrophenol	56	64	40 - 115	13	20	
2,4-Dimethylphenol	40	48	30 - 110	18	20	
Bis(2-chloroethoxy)methane	52	60	45 - 105	14	20	
2,4-Dichlorophenol	54	63	50 - 105	15	20	
1,2,4-Trichlorobenzene	54	60	35 - 105	10	20	
Naphthalene	55	62	40 - 100	12	20	
4-Chloroaniline	35	40	15 - 110	14	20	
Hexachlorobutadiene	50	54	25 - 105	8	20	
4-Chloro-3-methylphenol	51	62	45 - 110	20	20	
2-Methylnaphthalene	53	60	45 - 105	12	20	
Hexachlorocyclopentadiene	32	35	10 - 120	NC	20	U
2,4,6-Trichlorophenol	56	64	50 - 115	13	20	
2,4,5-Trichlorophenol	57	67	50 - 110	16	20	
2-Chloronaphthalene	57	61	50 - 105	6	20	
2-Nitroaniline	56	65	50 - 115	15	20	
Dimethyl phthalate	58	70	25 - 125	18	20	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 560-3816

**Method: 8270C
Preparation: 3520C**

MS Lab Sample ID: 560-1385-1
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1325
Date Prepared: 08/08/2006 1230

Analysis Batch: 560-3854
Prep Batch: 560-3816

Instrument ID: Agilent GCMS [Method
Lab File ID: 08090607.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

MSD Lab Sample ID: 560-1385-1
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1353
Date Prepared: 08/08/2006 1230

Analysis Batch: 560-3854
Prep Batch: 560-3816

Instrument ID: Agilent GCMS [Method 827
Lab File ID: 08090608.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

Analyte	% Rec.		Limit	RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD					
Acenaphthylene	58	63	50 - 105	8	20		
2,6-Dinitrotoluene	59	70	50 - 115	17	20		
3-Nitroaniline	53	64	20 - 125	18	20		
Acenaphthene	58	63	45 - 110	9	20		
2,4-Dinitrophenol	70	87	15 - 140	22	20		F
4-Nitrophenol	67	80	20 - 120	18	20		
2,4-Dinitrotoluene	62	74	50 - 120	17	20		
Diethyl phthalate	62	73	40 - 120	17	20		
Fluorene	59	66	50 - 110	11	20		
4-Chlorophenyl phenyl ether	57	63	50 - 110	10	20		
4-Nitroaniline	61	73	35 - 120	19	20		
4,6-Dinitro-2-methylphenol	71	90	40 - 130	23	20		F
N-Nitrosodiphenylamine	48	59	50 - 110	19	20		
4-Bromophenyl phenyl ether	57	67	50 - 115	17	20		
Hexachlorobenzene	58	67	50 - 110	15	20		
Phenanthrene	61	72	50 - 115	16	20		
Anthracene	61	71	55 - 110	16	20		
Di-n-butyl phthalate	61	71	55 - 115	16	20		
Fluoranthene	61	70	55 - 115	14	20		
Pyrene	58	70	50 - 130	19	20		
Butyl benzyl phthalate	57	68	45 - 115	16	20		
Benzo[a]anthracene	59	70	55 - 110	17	20		
Chrysene	59	69	55 - 110	15	20		
Bis(2-ethylhexyl) phthalate	58	68	40 - 125	15	20		
Di-n-octyl phthalate	60	69	35 - 135	15	20		
Benzo[b]fluoranthene	61	71	45 - 120	15	20		
Benzo[k]fluoranthene	55	66	45 - 125	17	20		
Benzo[a]pyrene	57	68	55 - 110	17	20		
Indeno[1,2,3-cd]pyrene	61	72	45 - 125	15	20		

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 560-3816

Method: 8270C
Preparation: 3520C

MS Lab Sample ID: 560-1385-1 Analysis Batch: 560-3854
Client Matrix: Water Prep Batch: 560-3816
Dilution: 1.0
Date Analyzed: 08/09/2006 1325
Date Prepared: 08/08/2006 1230

Instrument ID: Agilent GCMS [Method
Lab File ID: 08090607.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

MSD Lab Sample ID: 560-1385-1 Analysis Batch: 560-3854
Client Matrix: Water Prep Batch: 560-3816
Dilution: 1.0
Date Analyzed: 08/09/2006 1353
Date Prepared: 08/08/2006 1230

Instrument ID: Agilent GCMS [Method 827
Lab File ID: 08090608.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 1 mL
Injection Volume:

Analyte	% Rec.		Limit	RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD					
Dibenz(a,h)anthracene	60	71	40 - 125	16	20		
Benzo[g,h,i]perylene	60	69	40 - 125	15	20		
3,3'-Dichlorobenzidine	30	39	20 - 110	25	20		F
Pentachlorophenol	66	79	40 - 115	18	20		
N-Nitrosodimethylamine	46	50	25 - 110	8	20		
Benzoic acid	41	50	10 - 121	20	20	J	
Surrogate	MS % Rec		MSD % Rec		Acceptance Limits		
2-Fluorophenol	47		53		10 - 120		
Phenol-d5	50		56		12 - 120		
Nitrobenzene-d5	54		59		30 - 120		
2-Fluorobiphenyl	50		55		26 - 120		
2,4,6-Tribromophenol	57		70		25 - 120		
Terphenyl-d14	28		35		10 - 120		

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3832

Lab Sample ID: MB 560-3832/1-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/14/2006 1826
Date Prepared: 08/08/2006 1700

Analysis Batch: 560-3996
Prep Batch: 560-3832
Units: ug/L

Method: 8081A Preparation: 3520C

Instrument ID: Agilent GC [Method 8081]
Lab File ID: 08140637.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 10 mL
Injection Volume:
Column ID: PRIMARY

Analyte	Result	Qual	MDL	RL
alpha-BHC	0.0056	U	0.0056	0.050
beta-BHC	0.0056	U	0.0056	0.050
delta-BHC	0.0025	U	0.0025	0.050
Heptachlor	0.0059	U	0.0059	0.050
Aldrin	0.0025	U	0.0025	0.050
Heptachlor epoxide	0.0028	U	0.0028	0.050
4,4'-DDE	0.0026	U	0.0026	0.050
Endosulfan I	0.0089	U	0.0089	0.050
Dieldrin	0.0083	U	0.0083	0.050
Endrin	0.0025	U	0.0025	0.050
4,4'-DDD	0.0029	U	0.0029	0.050
Endosulfan II	0.0035	U	0.0035	0.050
4,4'-DDT	0.0034	U	0.0034	0.050
Methoxychlor	0.023	U	0.023	0.050
Endosulfan sulfate	0.0039	U	0.0039	0.050
Endrin ketone	0.0073	U	0.0073	0.050
Chlordane (technical)	0.050	U	0.050	0.50
Toxaphene	0.50	U	0.50	5.0
gamma-BHC (Lindane)	0.0027	U	0.0027	0.050
Surrogate	% Rec	Acceptance Limits		
Tetrachloro-m-xylene	73	57 - 127		
DCB Decachlorobiphenyl	63	10 - 152		

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Control Spike - Batch: 560-3832

Method: 8081A

Preparation: 3520C

Lab Sample ID: LCS 560-3832/2-B

Analysis Batch: 560-3996

Instrument ID: Agilent GC [Method 8081]

Client Matrix: Water

Prep Batch: 560-3832

Lab File ID: 08140639.D

Dilution: 1.0

Units: ug/L

Initial Weight/Volume: 1000 mL

Date Analyzed: 08/14/2006 1850

Final Weight/Volume: 10 mL

Date Prepared: 08/08/2006 1700

Injection Volume:

Column ID: PRIMARY

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
alpha-BHC	0.500	0.41	83	60 - 130	
beta-BHC	0.500	0.41	83	65 - 125	
delta-BHC	0.500	0.41	82	45 - 135	
Heptachlor	0.500	0.40	79	40 - 130	
Aldrin	0.500	0.40	81	25 - 140	
Heptachlor epoxide	0.500	0.42	84	60 - 130	
4,4'-DDE	0.500	0.44	87	35 - 140	
Endosulfan I	0.500	0.41	81	50 - 110	
Dieldrin	0.500	0.43	86	60 - 130	
Endrin	0.500	0.39	78	55 - 135	
4,4'-DDD	0.500	0.44	88	25 - 150	
Endosulfan II	0.500	0.42	84	30 - 130	
4,4'-DDT	0.500	0.46	92	45 - 140	
Methoxychlor	0.500	0.46	93	55 - 150	
Endosulfan sulfate	0.500	0.45	89	55 - 135	
Endrin ketone	0.500	0.47	95	75 - 125	
gamma-BHC (Lindane)	0.500	0.43	85	25 - 135	
Surrogate		% Rec		Acceptance Limits	
Tetrachloro-m-xylene		77		57 - 127	
DCB Decachlorobiphenyl		66		10 - 152	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 560-3832

MS Lab Sample ID: 560-1385-3
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/14/2006 2025
 Date Prepared: 08/08/2006 1700

MSD Lab Sample ID: 560-1385-3
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/14/2006 2048
 Date Prepared: 08/08/2006 1700

Analysis Batch: 560-3996
 Prep Batch: 560-3832

Analysis Batch: 560-3996
 Prep Batch: 560-3832

Method: 8081A
Preparation: 3520C

Instrument ID: Agilent GC [Method 8081]
 Lab File ID: 08140647.D
 Initial Weight/Volume: 1000 mL
 Final Weight/Volume: 10 mL
 Injection Volume:

Column ID: PRIMARY

Instrument ID: Agilent GC [Method 8081]
 Lab File ID: 08140649.D
 Initial Weight/Volume: 1000 mL
 Final Weight/Volume: 10 mL
 Injection Volume:

Column ID: PRIMARY

Analyte	% Rec.		Limit	RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD					
alpha-BHC	78	75	60 - 130	4	30		
beta-BHC	72	68	65 - 125	6	30		
delta-BHC	69	63	45 - 135	9	30		
Heptachlor	51	42	40 - 130	20	30		
Aldrin	45	37	25 - 140	20	30		
Heptachlor epoxide	63	54	60 - 130	15	30	F	
4,4'-DDE	38	30	35 - 140	22	30	F	
Endosulfan I	61	53	50 - 110	15	30		
Dieldrin	66	56	60 - 130	17	30	F	
Endrin	71	60	55 - 135	17	30		
4,4'-DDD	55	50	25 - 150	9	30		
Endosulfan II	62	53	30 - 130	15	30		
4,4'-DDT	46	37	45 - 140	22	30	F	
Methoxychlor	62	49	55 - 150	23	30	F	
Endosulfan sulfate	79	68	55 - 135	15	30		
Endrin ketone	79	69	75 - 125	14	30	F	
gamma-BHC (Lindane)	77	73	25 - 135	5	30		
Surrogate	MS % Rec		MSD % Rec		Acceptance Limits		
Tetrachloro-m-xylene	63		60		57 - 127		
DCB Decachlorobiphenyl	28		24		10 - 152		

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Enact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3831

Method: 8082

Preparation: 3520C

Lab Sample ID: MB 560-3831/1-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/15/2006 1904
Date Prepared: 08/08/2006 1700

Analysis Batch: 560-4015
Prep Batch: 560-3831
Units: ug/L

Instrument ID: Hewlett Packard GC [Meth]
Lab File ID: 08150632.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 10 mL
Injection Volume:
Column ID: PRIMARY

Analyte	Result	Qual	MDL	RL
Aroclor 1016	0.17	U	0.17	0.50
Aroclor 1221	0.17	U	0.17	0.50
Aroclor 1232	0.17	U	0.17	0.50
Aroclor 1242	0.17	U	0.17	0.50
Aroclor 1248	0.17	U	0.17	0.50
Aroclor 1254	0.17	U	0.17	0.50
Aroclor 1260	0.17	U	0.17	0.50

Surrogate	% Rec	Acceptance Limits
Tetrachloro-m-xylene	96	25 - 140
DCB Decachlorobiphenyl	82	42 - 133

Method Blank - Batch: 560-3831

Method: 8082

Preparation: 3520C

Lab Sample ID: MB 560-3831/1-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/16/2006 1630
Date Prepared: 08/08/2006 1700

Analysis Batch: 560-4015
Prep Batch: 560-3831
Units: ug/L

Instrument ID: Hewlett Packard GC [Meth]
Lab File ID: 08160607.D
Initial Weight/Volume: 1000 mL
Final Weight/Volume: 10 mL
Injection Volume:
Column ID: PRIMARY

Analyte	Result	Qual	MDL	RL
Aroclor 1016	0.17	U	0.17	0.50
Aroclor 1221	0.17	U	0.17	0.50
Aroclor 1232	0.17	U	0.17	0.50
Aroclor 1242	0.17	U	0.17	0.50
Aroclor 1248	0.17	U	0.17	0.50
Aroclor 1254	0.17	U	0.17	0.50
Aroclor 1260	0.17	U	0.17	0.50

Surrogate	% Rec	Acceptance Limits
Tetrachloro-m-xylene	89	25 - 140
DCB Decachlorobiphenyl	70	42 - 133

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Control Spike - Batch: 560-3831

Method: 8082

Preparation: 3520C

Lab Sample ID: LCS 560-3831/2-A

Analysis Batch: 560-4015

Instrument ID: Hewlett Packard GC [Meth]

Client Matrix: Water

Prep Batch: 560-3831

Lab File ID: 08150633.D

Dilution: 1.0

Units: ug/L

Initial Weight/Volume: 1000 mL

Date Analyzed: 08/15/2006 1922

Final Weight/Volume: 10 mL

Date Prepared: 08/08/2006 1700

Injection Volume:

Column ID: PRIMARY

Analyte	Spike Amount	Result	% Rec.	Limit	Qual					
Aroclor 1016	10.0	10.6	106	50 - 135						
Aroclor 1260	10.0	9.63	96	50 - 135						
Surrogate	% Rec									
Tetrachloro-m-xylene	100									
DCB Decachlorobiphenyl	83									
Acceptance Limits										
25 - 140										
42 - 133										

Matrix Spike/

Matrix Spike Duplicate Recovery Report - Batch: 560-3831

Method: 8082

Preparation: 3520C

MS Lab Sample ID: 560-1385-2

Analysis Batch: 560-4015

Instrument ID: Hewlett Packard GC [Met

Client Matrix: Water

Prep Batch: 560-3831

Lab File ID: 08150636.D

Dilution: 1.0

Initial Weight/Volume: 1000 mL

Date Analyzed: 08/15/2006 2013

Final Weight/Volume: 10 mL

Date Prepared: 08/08/2006 1700

Injection Volume:

Column ID: PRIMARY

MSD Lab Sample ID: 560-1385-2

Analysis Batch: 560-4015

Instrument ID: Hewlett Packard GC [Meth]

Client Matrix: Water

Prep Batch: 560-3831

Lab File ID: 08150637.D

Dilution: 1.0

Initial Weight/Volume: 1000 mL

Date Analyzed: 08/15/2006 2030

Final Weight/Volume: 10 mL

Date Prepared: 08/08/2006 1700

Injection Volume:

Column ID: PRIMARY

Analyte	<u>% Rec.</u>		Limit	RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD					
Aroclor 1016	84	83	50 - 135	0	30		
Aroclor 1260	33	37	50 - 135	12	30	F	F
Surrogate	<u>MS % Rec</u>					<u>Acceptance Limits</u>	
Tetrachloro-m-xylene	81		83			25 - 140	
DCB Decachlorobiphenyl	24	X	33	X		42 - 133	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Enact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3757

Lab Sample ID: MB 560-3757/1-A
Client Matrix: Water
Dilution: 10
Date Analyzed: 08/09/2006 0050
Date Prepared: 08/07/2006 1103

Analysis Batch: 560-3885
Prep Batch: 560-3757
Units: ug/L

Method: 6020
Preparation: 3010A

Instrument ID: Agilent ICPMS
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Analyte

Analyte	Result	Qual	MDL	RL
Ag	1.0	U	1.0	5.0
As	1.0	U	1.0	5.0
Ba	12	J	1.0	50
Cd	1.0	U	1.0	5.0
Cr	1.1	U	1.1	20
Ni	3.3	J	1.0	5.0
Pb	1.0	U	1.0	5.0
Se	1.0	U	1.0	5.0
Zn	50	U	50	100

Method Blank - Batch: 560-3757

Lab Sample ID: MB 560-3757/25-A
Client Matrix: Water
Dilution: 10
Date Analyzed: 08/09/2006 0457
Date Prepared: 08/07/2006 1103

Analysis Batch: 560-3885
Prep Batch: 560-3757
Units: ug/L

Method: 6020
Preparation: 3010A
Dissolved

Instrument ID: Agilent ICPMS
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Analyte

Analyte	Result	Qual	MDL	RL
Ag	1.0	U	1.0	5.0
As	1.0	U	1.0	5.0
Ba	2.8	J	1.0	50
Cd	1.0	U	1.0	5.0
Cr	1.1	U	1.1	20
Ni	1.1	J	1.0	5.0
Pb	1.0	U	1.0	5.0
Se	1.0	U	1.0	5.0
Zn	50	U	50	100

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Lab Control Spike - Batch: 560-3757

Method: 6020
Preparation: 3010A

Lab Sample ID: LCS 560-3757/2-A
Client Matrix: Water
Dilution: 10
Date Analyzed: 08/09/2006 0056
Date Prepared: 08/07/2006 1103

Analysis Batch: 560-3885
Prep Batch: 560-3757
Units: ug/L

Instrument ID: Agilent ICPMS
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Ag	200	190	97	80 - 120	
As	400	430	108	80 - 120	
Ba	400	430	106	80 - 120	
Cd	200	220	109	80 - 120	
Cr	400	420	106	80 - 120	
Ni	400	430	108	80 - 120	
Pb	200	200	102	80 - 120	
Se	400	490	124	80 - 120	
Zn	400	440	109	80 - 120	

Lab Control Spike - Batch: 560-3757

Method: 6020
Preparation: 3010A
Dissolved

Lab Sample ID: LCS 560-3757/26-A
Client Matrix: Water
Dilution: 10
Date Analyzed: 08/09/2006 0503
Date Prepared: 08/07/2006 1103

Analysis Batch: 560-3885
Prep Batch: 560-3757
Units: ug/L

Instrument ID: Agilent ICPMS
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Ag	200	210	106	80 - 120	
As	400	500	125	80 - 120	*
Ba	400	490	121	80 - 120	*
Cd	200	250	125	80 - 120	*
Cr	400	500	124	80 - 120	*
Ni	400	510	128	80 - 120	*
Pb	200	230	115	80 - 120	
Se	400	550	138	80 - 120	*
Zn	400	510	128	80 - 120	*

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 560-3757

MS Lab Sample ID: 560-1385-1 Analysis Batch: 560-3885
Client Matrix: Water Prep Batch: 560-3757
Dilution: 10
Date Analyzed: 08/09/2006 0515
Date Prepared: 08/07/2006 1103

MSD Lab Sample ID: 560-1385-1 Analysis Batch: 560-3885
Client Matrix: Water Prep Batch: 560-3757
Dilution: 10
Date Analyzed: 08/09/2006 0521
Date Prepared: 08/07/2006 1103

Method: 6020
Preparation: 3010A
Dissolved

Instrument ID: Agilent ICPMS
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Instrument ID: Agilent ICPMS
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Analyte	% Rec.		Limit	RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD					
Ag	87	94	75 - 125	8	20		
As	110	110	75 - 125	0	20		
Ba	89	100	75 - 125	8	20		
Cd	104	110	75 - 125	6	20		
Cr	104	104	75 - 125	0	20		
Ni	105	106	75 - 125	1	20		
Pb	100	104	75 - 125	3	20		
Se	122	116	75 - 125	5	20		
Zn	108	109	75 - 125	1	20		

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3833

Lab Sample ID: MB 560-3833/19-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1607
Date Prepared: 08/09/2006 1130

Analysis Batch: 560-3844
Prep Batch: 560-3833
Units: mg/L

Method: 7470A
Preparation: 7470A
Dissolved

Instrument ID: Mercury Analyzer Leeman
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Analyte	Result	Qual	MDL	RL
Hg	0.00013	U	0.00013	0.0020

Lab Control Spike - Batch: 560-3833

Lab Sample ID: LCS 560-3833/18-A
Client Matrix: Water
Dilution: 1.0
Date Analyzed: 08/09/2006 1614
Date Prepared: 08/09/2006 1130

Analysis Batch: 560-3844
Prep Batch: 560-3833
Units: mg/L

Method: 7470A
Preparation: 7470A
Dissolved

Instrument ID: Mercury Analyzer Leeman
Lab File ID: N/A
Initial Weight/Volume: 50 mL
Final Weight/Volume: 50 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Hg	0.00500	0.00507	101	80 - 120	

Calculations are performed before rounding to avoid round-off errors in calculated results.

Quality Control Results

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Method Blank - Batch: 560-3912

Lab Sample ID: MB 560-3912/4-A
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/11/2006 1532
 Date Prepared: 08/11/2006 1105

Analysis Batch: 560-3921
 Prep Batch: 560-3912
 Units: mg/L

Method: 7470A
Preparation: 7470A
Dissolved

Instrument ID: Mercury Analyzer Leeman
 Lab File ID: N/A
 Initial Weight/Volume: 50 mL
 Final Weight/Volume: 50 mL

Analyte	Result	Qual	MDL	RL
Hg	0.00013	U	0.00013	0.0020

Lab Control Spike - Batch: 560-3912

Lab Sample ID: LCS 560-3912/3-A
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/11/2006 1535
 Date Prepared: 08/11/2006 1105

Analysis Batch: 560-3921
 Prep Batch: 560-3912
 Units: mg/L

Method: 7470A
Preparation: 7470A

Instrument ID: Mercury Analyzer Leeman
 Lab File ID: N/A
 Initial Weight/Volume: 50 mL
 Final Weight/Volume: 50 mL

Analyte	Spike Amount	Result	% Rec.	Limit	Qual
Hg	0.00500	0.0047	94	80 - 120	

Matrix Spike/ Matrix Spike Duplicate Recovery Report - Batch: 560-3912

MS Lab Sample ID: 560-1385-10
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/11/2006 1541
 Date Prepared: 08/11/2006 1105

Analysis Batch: 560-3921
 Prep Batch: 560-3912

Method: 7470A
Preparation: 7470A
Dissolved

Instrument ID: Mercury Analyzer Leeman
 Lab File ID: N/A
 Initial Weight/Volume: 50 mL
 Final Weight/Volume: 50 mL

MSD Lab Sample ID: 560-1385-10
 Client Matrix: Water
 Dilution: 1.0
 Date Analyzed: 08/11/2006 1544
 Date Prepared: 08/11/2006 1105

Analysis Batch: 560-3921
 Prep Batch: 560-3912

Instrument ID: Mercury Analyzer Leeman
 Lab File ID: N/A
 Initial Weight/Volume: 50 mL
 Final Weight/Volume: 50 mL

Analyte	% Rec.		Limit	RPD	RPD Limit	MS Qual	MSD Qual
	MS	MSD					
Hg	80	83	80 - 120	3	20		

Calculations are performed before rounding to avoid round-off errors in calculated results.

SEVERN
TRENT

STL

No. 38281

CHAIN OF CUSTODY RECORD

CUSTOMER INFORMATION		PROJECT INFORMATION					NUMBER OF CONTAINERS	ANALYSIS/METHOD REQUEST							
COMPANY:	Entact	PROJECT NAME/NUMBER: Sheridan /D1631						VOCs	SVOCs	8260-3	8270C	METALS	Pesticides	PCBs	Dust/PAHs
SEND REPORT TO:	Liz Scaggs	BILLING INFORMATION													
ADDRESS:	3129 Bass Pro Dr Grapevine, TX 76051	BILL TO:	Sevme												
PHONE:	972 580 1323	ADDRESS:													
FAX:	972 550 7464	PHONE:													
SAMPLE NO.	SAMPLE DESCRIPTION	SAMPLE DATE	SAMPLE TIME	SAMPLE MATRIX	CONTAINER	PRESERV.									
1	R2-A	8-2-06	1123	H ₂ O	1000 ML STL	ice HNO ₃ ice		II	✓	✓	✓	✓	✓	✓	✓
2	R2-B		1147				II	✓	✓	✓	✓	✓	✓		
3	R2-C		1215				II	✓	✓	✓	✓	✓	✓		
4	R2-D		1241				II	✓	✓	✓	✓	✓	✓		
5	R1-A		1621				II	✓	✓	✓	✓	✓	✓		
6	R1-B		1643				II	✓	✓	✓	✓	✓	✓		
7	R1-C		1709				II	✓	✓	✓	✓	✓	✓		
8	R1-D		↓ 1729	↓	↓	↓	II	✓	✓	✓	✓	✓	✓		
SAMPLER: D. McGough / R. Smith		SHIPMENT METHOD: FedEx					AIRBILL NO.:								
REQUIRED TURNAROUND: <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HOURS <input type="checkbox"/> 48 HOURS <input type="checkbox"/> 72 HOURS <input type="checkbox"/> 5 DAYS <input type="checkbox"/> 10 DAYS <input checked="" type="checkbox"/> ROUTINE <input type="checkbox"/> OTHER															
1. RELINQUISHED BY: SIGNATURE: <i>JK Robyn Smith</i>		DATE: 8-3-06	2. RELINQUISHED BY: SIGNATURE: <i>Ted ey</i>		DATE	3. RELINQUISHED BY: SIGNATURE:		DATE							
PRINTED NAME/COMPANY: Entact		TIME: 1730	PRINTED NAME/COMPANY: <i>J. Robyn Smith</i>		TIME	PRINTED NAME/COMPANY:		TIME							
1. RECEIVED BY: SIGNATURE: <i>Ted ey</i>		DATE	2. RECEIVED BY: SIGNATURE: <i>J. Robyn Smith</i>		DATE	3. RECEIVED BY: SIGNATURE:		DATE							
PRINTED NAME/COMPANY: STL		TIME: 10:56A	PRINTED NAME/COMPANY: <i>STL</i>		TIME	PRINTED NAME/COMPANY:		TIME							

SEVERN TRENT LABORATORIES, INC. *LP*

1733 N. Padre Island Drive
 Corpus Christi, TX 78408
 Phone: (361) 289-2673 / Fax: (361) 289-2471

STL8222-560 (12/02)

SEVERN
TRENT

STL

2.2, 2.0, 1.0, 1.2, 1.0, 1.3, 0.7,
0.4, 2.2, 0.6, 1.4, 0.4, 0.2, 0.5 No. 38282

CHAIN OF CUSTODY RECORD

CUSTOMER INFORMATION		PROJECT INFORMATION		NUMBER OF CONTAINERS	ANALYSIS/METHOD REQUEST	LAB JOB NO.
COMPANY:	Entact	PROJECT NAME/NUMBER:	Sheridan 1D1631			
SEND REPORT TO:	Liz Scaggs	BILLING INFORMATION				
ADDRESS:	3129 Bass Pro Dr Grapevine, TX 76051	BILL TO:	Game			
ADDRESS:		ADDRESS:				
PHONE:	972 580 1323	PHONE:				
FAX:	972 550 7464	FAX:	PO NO:			
SAMPLE NO.	SAMPLE DESCRIPTION	SAMPLE DATE	SAMPLE TIME			
1	MW6	8-3-06	759	H ₂ O	1000mL 40mL	HNO ₃ HCl ICE
2	MW31		920			
3	MW34		1040			
4	MW35		1155			
5	MW37		1253			
6	MW39		1358	↓	↓	↓
SAMPLER:	D. McGough / R. Smith	SHIPMENT METHOD:	FedEx		AIRBILL NO.:	
REQUIRED TURNAROUND: <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HOURS <input type="checkbox"/> 48 HOURS <input type="checkbox"/> 72 HOURS <input type="checkbox"/> 5 DAYS <input type="checkbox"/> 10 DAYS <input checked="" type="checkbox"/> ROUTINE <input type="checkbox"/> OTHER						
1. RELINQUISHED BY:	DATE	2. RELINQUISHED BY:	DATE	3. RELINQUISHED BY:	DATE	
SIGNATURE: K. Robyn Smith	8-3-06	SIGNATURE: Jd ex		SIGNATURE:		
PRINTED NAME/COMPANY: Entact	TIME: 1730	PRINTED NAME/COMPANY:	TIME	PRINTED NAME/COMPANY:	TIME	
1. RECEIVED BY:	DATE	2. RECEIVED BY:	DATE	3. RECEIVED BY:	DATE	
SIGNATURE: Jd ex		SIGNATURE: P. D. WILSON	8/4/06	SIGNATURE:		
PRINTED NAME/COMPANY:	TIME	PRINTED NAME/COMPANY:	TIME	PRINTED NAME/COMPANY:	TIME	

SEVERN TRENT LABORATORIES, INC. LP

1733 N. Padre Island Drive
Corpus Christi, TX 78408
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STL8222-560 (12/02)

LOGIN SAMPLE RECEIPT CHECK LIST

Client: Entact Environmental Services, LLC

Job Number: 560-1385-1

Login Number: 1385

Question	T/F/NA	Comment
Radioactivity either was not measured or, if measured, is at or below background	NA	
The cooler's custody seal, if present, is intact.	True	
The cooler or samples do not appear to have been compromised or tampered with.	True	
Samples were received on ice.	True	
Cooler Temperature is acceptable.	True	
Cooler Temperature is recorded.	True	
COC is present.	True	
COC is filled out in ink and legible.	True	
COC is filled out with all pertinent information.	True	
There are no discrepancies between the sample IDs on the containers and the COC.	True	Note: SX R2-A & R2-B HAD CONTAINERS W/RIVER 2 SET A & RIVER 2 SET B ON LABEL
Samples are received within Holding Time.	True	
Sample containers have legible labels.	True	
Containers are not broken or leaking.	True	
Sample collection date/times are provided.	True	
Appropriate sample containers are used.	True	
Sample bottles are completely filled.	True	
There is sufficient vol. for all requested analyses, incl. any requested MS/MSDs	True	
VOA sample vials do not have headspace or bubble is <6mm (1/4") in diameter.	True	
If necessary, staff have been informed of any short hold time or quick TAT needs	True	
Multiphasic samples are not present.	True	
Samples do not require splitting or compositing.	True	

SEVERN
TRENT

STL

No. 38281

CHAIN OF CUSTODY RECORD

CUSTOMER INFORMATION		PROJECT INFORMATION		NUMBER OF CONTAINERS	ANALYSIS/METHOD REQUEST								
COMPANY:	Entact	PROJECT NAME/NUMBER:	Sheridan / 101631			VOCS	SVOCs	8260 B	8270C	8010B	17470A	PCBs	8081A
SEND REPORT TO:	Liz Scagg's	BILL TO:	same										
ADDRESS:	3129 Bass Pro Dr Grapvine, TX 76051	ADDRESS:											
PHONE:	972 580 1323	PHONE:											
FAX:	972 550 7464	FAX:											
PO NO:		PO NO:											
SAMPLE NO.	SAMPLE DESCRIPTION	SAMPLE DATE	SAMPLE TIME			SAMPLE MATRIX	CONTAINER	PRESERV					
1	R2-A	8-2-06	1123			H2O	1000 ML TOTAL	1000 ML H2O & ICE	11	✓	✓	✓	✓
2	R2-B		1147				11	✓	✓	✓	✓		
3	R2-C		1215				11	✓	✓	✓	✓		
4	R2-D		1241				11	✓	✓	✓	✓		
5	R1-A		1621				11	✓	✓	✓	✓		
6	R1-B		1643				11	✓	✓	✓	✓		
7	R1-C		1709				11	✓	✓	✓	✓		
8	R1-D		1729	↓	↓	↓	11	✓	✓	✓	✓		
SAMPLER:	D. McGough / R. Smith	SHIPMENT METHOD:		FedEx		AIRBILL NO.:							
REQUIRED TURNAROUND* <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HOURS <input type="checkbox"/> 48 HOURS <input type="checkbox"/> 72 HOURS <input type="checkbox"/> 5 DAYS <input type="checkbox"/> 10 DAYS <input checked="" type="checkbox"/> ROUTINE <input type="checkbox"/> OTHER													
1. RELINQUISHED BY:	DATE	2. RELINQUISHED BY:	DATE	3. RELINQUISHED BY:	DATE								
SIGNATURE: <i>DR Robyn Smith</i>	8-3-06	SIGNATURE: <i>Ted ey</i>		SIGNATURE:									
PRINTED NAME/COMPANY: Entact	TIME 1730	PRINTED NAME/COMPANY: <i>J. Plourneau</i>	TIME	PRINTED NAME/COMPANY:	TIME								
1. RECEIVED BY:	DATE	2. RECEIVED BY:	DATE	3. RECEIVED BY:	DATE								
SIGNATURE: <i>Ted ey</i>		SIGNATURE: <i>J. Plourneau</i>	8/4/06	SIGNATURE:									
PRINTED NAME/COMPANY:	TIME	PRINTED NAME/COMPANY: <i>STL</i>	TIME 10:50	PRINTED NAME/COMPANY:	TIME								

SEVERN TRENT LABORATORIES, INC.

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 Corpus Christi, TX 78408
 Phone: (361) 289-2673 / Fax: (361) 289-2471

STL8222-560 (12/02)

SEVERN TRENT

STL

2.2, 2.0, 1.0, 1.4, 1.0, 1.0, 0.4, 0.2, 0.5 No. 38282
0.4, 2.2, 1.4, 1.4, 0.4, 0.2, 0.5
CHAIN OF CUSTODY RECORD

CHAIN OF CUSTODY RECORD

CUSTOMER INFORMATION		PROJECT INFORMATION		NUMBER OF CONTAINERS	ANALYSIS METHOD REQUEST											
COMPANY:	Entact	PROJECT NAME/NUMBER:	Sheridan 1D1631													
SEND REPORT TO:	Liz Scaggs	BILLING INFORMATION														
ADDRESS:	3129 Bass Pro DR Grapevine, TX 76051	BILL TO:	Same													
PHONE:	972 580 1323	PHONE:														
FAX:	972 550 7464	FAX:														
SAMPLE NO.	SAMPLE DESCRIPTION	SAMPLE DATE	SAMPLE TIME			SAMPLE MATRIX	CONTAINER	PRESERV	11	✓	✓	✓	✓	✓	✓	REMARKS/PRECAUTIONS
1	MW6	8.3.06	759			H ₂ O	1000ML YONAL	HNO ₃ HCL ICE	11	✓	✓	✓	✓	✓	✓	two metal samples submitted
2	MW31		920				11	✓	✓	✓	✓	✓	✓	one unfiltered		
3	MW34		1040				11	✓	✓	✓	✓	✓	✓	& one field filtered with		
4	MW35		1155				11	✓	✓	✓	✓	✓	✓	45 micron filter		
5	MW37		1253				11	✓	✓	✓	✓	✓	✓			
6	MW39	↓	1358	↓	↓	↓	11	✓	✓	✓	✓	✓	✓			
SAMPLER:	D. McGough / R. Smith	SHIPMENT METHOD:		FedEx				AIRBILL NO.:								
REQUIRED TURNAROUND* <input type="checkbox"/> SAME DAY <input type="checkbox"/> 24 HOURS <input type="checkbox"/> 48 HOURS <input type="checkbox"/> 72 HOURS <input type="checkbox"/> 5 DAYS <input type="checkbox"/> 10 DAYS <input checked="" type="checkbox"/> ROUTINE <input type="checkbox"/> OTHER																
1. RELINQUISHED BY:	DATE	2. RELINQUISHED BY:	DATE	3. RELINQUISHED BY:	DATE											
SIGNATURE: K. Robyn Smith	8.3.06	SIGNATURE: Ted ey		SIGNATURE:												
PRINTED NAME/COMPANY: Entact	TIME: 1730	PRINTED NAME/COMPANY:	TIME	PRINTED NAME/COMPANY:	TIME											
1. RECEIVED BY:	DATE	2. RECEIVED BY:	DATE	3. RECEIVED BY:	DATE											
SIGNATURE: Ted ey		SIGNATURE: P. DUMMAM	8/4/06	SIGNATURE:												
PRINTED NAME/COMPANY:	TIME	PRINTED NAME/COMPANY:	TIME	PRINTED NAME/COMPANY:	TIME											

***RUSH TURNAROUND MAY REQUIRE SURCHARGE**

SEVERN TRENT LABORATORIES, INC.

1733 N. Padre Island Dr.
Corpus Christi, TX 7840

Phone: (361) 289-2673 / Fax: (361) 289-2471

STL8222-560 (12/02)



ENTACT

Appendix

B

Appendix B

Appendix B
Statistical Calculations

Compound	Date	Benzene	Tetrachloroethylene	Trans-1,2-Dichloroethylene	Trichloroethylene	Total Arsenic
Alternate Concentration Limit		26	41	26	26	260
Trigger for RAP Preparation		4	6	4	4	40
Trigger for Increased Monitoring		1	2	1	1	10
R1-A ADJ	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0047
R1-B ADJ	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0045
R1-C ADJ	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0041
R1-D ADJ	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0054
R2-A ¹ DOWN	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0051
R2-B DOWN	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0058
R2-C DOWN	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0043
R2-D DOWN	08/02/06	0.0002	0.0002	0.0002	0.00032	0.0046

Note - all concentrations in mg/L

1 - Downgradient Brazos River Sample

J - Result is less than the RL but greater than or equal to the MDL and the concentration is an approximate value

DOWNSTREAM BACKGROUND WATER QUALITY

	Benzene	PCE	trans-1,2-DCE	TCE	Arsenic
Mean	0.0002	0.0002	0.0002	0.00032	0.00495
Variance	0.0002	0.0002	0.0002	0.00032	3.225E-07

ADJACENT BACKGROUND WATER QUALITY

	Benzene	PCE	trans-1,2-DCE	TCE	Arsenic
Mean	0.0002	0.0002	0.0002	0.00032	0.004675
Variance	0.0002	0.0002	0.0002	0.00032	2.2188E-07

NORMALITY DISTRIBUTION BY GEARY'S PROCEDURE

ALL	Benzene	PCE	trans-1,2-DCE	TCE	Arsenic
Mean					0.0048125
SSS					2.329E-06
SAD					3.725E-03
The Test					8.630E-01
Significance					0.86755296

DUNNETT'S PROCEDURE FOR DOWNSTREAM ARSENIC

	Adjacent	Downstream
Ex	0.0187	0.0198
xi	0.004675	0.00495
xi-xo		-0.000275
Ex2	0.00008831	0.0000993
Si2	0.00000030	0.00000043
Ti		